

## Inrobin

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## Experiment two:

### Detecting Spectral Content

Set the notation of this experiment as follows:

$\Omega$  is the set of  $M_{true}$  vectors of parameters where each vector is  $J$  dimensional

$$\Omega = \left\{ \Psi^1, \dots, \Psi^{M_{true}} \right\}$$

Let  $\Psi^j = [\Psi_1^j, \dots, \Psi_J^j]$  which is the  $j$ th element of  $\Omega$  set

$\Psi_h^j$  is  $h$ th element of  $\Psi^j$

$k(\cdot, \cdot; \Psi)$  a kernel function from parameterized by the parameter  $\Psi$ ,  $k : \mathcal{R} \times \mathcal{R} \rightarrow \mathcal{R}$

Let  $\mathbf{t}$  be a  $1 \times N$  vector of real numbers

$$\mathbf{K} = k(\mathbf{t}, \mathbf{t}; \Psi = \Psi)$$

$$j \in \left\{ 1, \dots, M_{true} \right\}, m \in \left\{ 1, \dots, M \right\}, h \in \left\{ 1, \dots, J \right\}$$

$n_{j,m}$  number of IMFs

#### Algorithm 1: Algorithm

**Input:** Define  $k, J, \Omega, \mathbf{t}, M_{true}$

Set  $m, j, J$

1. Evaluate each Gram Matrix  $\mathbf{K}_h^{(j)} = k(\mathbf{t}, \mathbf{t}; \Psi_h^{(j)})$  parametrized by the  $j$ th parameter from  $\Omega$ . There are  $J$  Gram Matrices since  $h = 1, \dots, J$
2. Simulate  $\mathbf{y}_{j,h}^{(m)} \sim \mathcal{N}(0, \mathbf{K}_h^{(j)})$  an  $N$  dimensional vector for each Gram Matrix  $\mathbf{K}_h^{(j)}$ , therefore we have  $J$  one dimensional vectors of length  $N$ .
3. Compute  $\mathbf{x}_j^{(m)} = \sum_{h=1}^J \mathbf{y}_{j,h}^{(m)}$
4. Fit a spline through each  $\mathbf{x}_j^{(m)}$  denoted as  $\hat{\mathbf{x}}_j^{(m)}$ .
5. Apply the EMD to  $\hat{\mathbf{x}}_j^{(m)}$  to get the IMFs decomposition and collect all the IMFs generated up to the stopping criterion chosen denoted as  $\gamma_{j,1}^{(m)}, \gamma_{j,2}^{(m)}, \dots, \gamma_{j,v_{j,m}}^{(m)}$ . For each  $j$  and for each  $m$  we might have different number of IMFs denote by  $v_{j,m}$ .
6. Compute the Instantaneous Frequency of each IMF denoted as  $f_{j,1}^{(m)}, f_{j,2}^{(m)}, \dots, f_{j,v_{j,m}}^{(m)}$ .
7. Compare the frequencies with Spectral Component of the kernels

**Algorithm 2:** Algorithm

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Input: Define  $k, J, \Omega, \mathbf{t}, M_{true}$ 
for ( $m = 1 : M$ ) {
    for ( $h = 1 : J$ ) {
        Evaluate each Gram Matrix  $\mathbf{K}_h^{(j)} = k(\mathbf{t}, \mathbf{t}; \Psi_h^{(j)})$  parametrized by the  $j$ th parameter from  $\Omega$ . There are  $J$  Gram Matrices since  $h = 1, \dots, J$ 
        Simulate  $\mathbf{y}_{j,h}^{(m)} \sim \mathcal{N}(0, \mathbf{K}_h^{(j)})$  an  $N$  dimensional vector for each Gram Matrix  $\mathbf{K}_h^{(j)}$ , therefore we have  $J$  one dimensional vectors of length  $N$ .
        Compute  $\mathbf{x}_j^{(m)} = \sum_{h=1}^J \mathbf{y}_{j,h}^{(m)}$ 
        Fit a spline through each  $\mathbf{x}_j^{(m)}$  denoted as  $\hat{\mathbf{x}}_j^{(m)}$ .
        Apply the EMD to  $\hat{\mathbf{x}}_j^{(m)}$  to get the IMFs decomposition and collect all the IMFs generated up to the stopping criterion chosen denoted as  $\gamma_{j,1}^{(m)}, \gamma_{j,2}^{(m)}, \dots, \gamma_{j,v_{j,m}}^{(m)}$ .
        For each  $j$  and for each  $m$  we might have different number of IMFs denote by  $v_{j,m}$ .
        Compute the Instantaneous Frequency of each IMF denoted as  $f_{j,1}^{(m)}, f_{j,2}^{(m)}, \dots, f_{j,v_{j,m}}^{(m)}$ .
    }
    Compare the frequencies with Spectral Component of the kernels

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Example: Periodic Kernel:  $\Omega = \{(l_1, p_1), (l_1, p_2), (l_1, p_3), (l_2, p_1), (l_2, p_2), (l_2, p_3)\}$ ,  $j = 2$ ,  $h = 1 : 6$ ,  $J = 6$ ,  $\Psi^j = [\Psi_1^j, \dots, \Psi_6^j]$  and  $\Psi_1^j = \{l_1, p_1\}, \dots, \Psi_6^j = \{l_2, p_3\}$ .

### Stopping Criteria for the sifting procedure

- **Cauchy-Type Convergence (SD).** [?] proposed such criterion to limit amplitude and frequency modulations through a certain threshold of the standard deviation of two consecutive sifting results.

$$SD = \sum_{t=0}^T \left[ \frac{|(h_{1(k-1)}(t) - h_{1k}(t))|^2}{h_{1(k-1)}^2(t)} \right] \quad (1)$$

- **Mean Fluctuations Threshold.** [?] proposed three different thresholds allowing for small fluctuations of the mean and locally large excursions. The thresholds are  $\alpha$ ,  $\theta_1$  and  $\theta_2$ . For  $(1 - \alpha)$  data, the criterion keeps sifting if  $\sigma(t) < \theta_1$ , while for the remaining fraction if  $\sigma(t) < \theta_2$ . The quantities are defines as:

$$\sigma(t) := \left| \frac{m(t)}{a(t)} \right| \quad a(t) := \frac{S(e_1^S) - S(e_2^S)}{2} \quad m(t) := \frac{S(e_1^S) + S(e_2^S)}{2}$$

3

- **Energy Difference Tracking.** The assumption of this criterion is that the EMD provides IMFs and a residual mutually orthogonal. If  $c_1$  is extracted from the signal, the sum of its energy and those of the residual signal ( $E_{tot}$ ) is equal to the energy of the original signal ( $E_S$ ). If not, there must be a difference called energy error ( $E_{err}$ ). The sifting procedure stops when both  $E_{err}$  and the mean value are small enough.  $E_{err}$  for  $c_1$  is defined as:

$$E_{err} = E_{tot} - E_S = \int c_1^2(t) dt - \int S(t)c_1(t) dt \quad (2)$$

- **Orthogonality Criterion.** [?] underline the fact that any IMF, in this case  $c_1$  should satisfy:

$$\sum_{t=1}^N c_1(t)(x(t) - c_1(t)) = 0 \quad (3)$$

which states the orthogonality of the IMFs. They determine the following index and the sifting procedure stops once that it reaches a certain minimum value:

$$OC = \left| \sum_{t=1}^N \frac{m(t) S(t)}{m(t)(S(t) - m(t))} \right| \quad \text{where} \quad m(t) = \frac{S(e_1^S) + S(e_2^S)}{2} \quad (4)$$

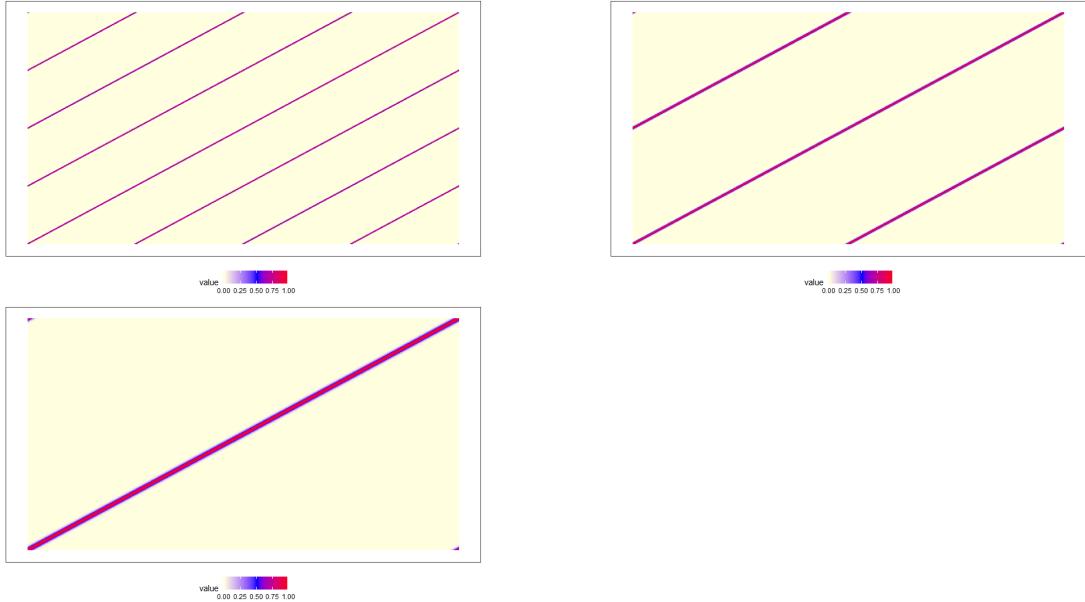


Figure 1: Heatmaps of Gram Matrices for the periodic Kernel with  $N = 1000, l = 0.05, p = 0.25, 0.5, 1$

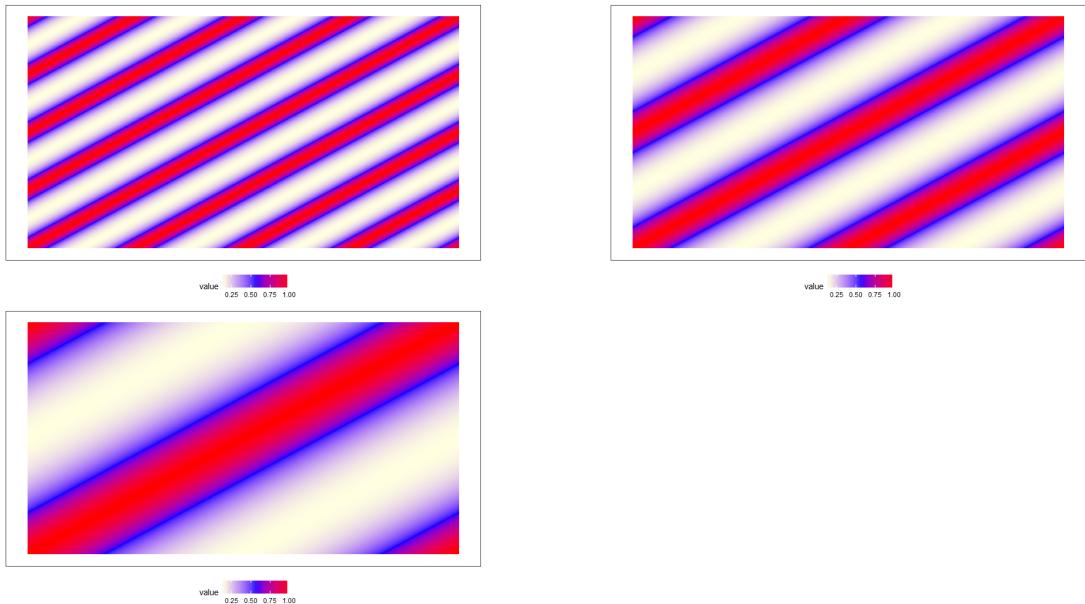


Figure 2: Heatmaps of Gram Matrices for the periodic Kernel with  $N = 1000, l = 1, p = 0.25, 0.5, 1$

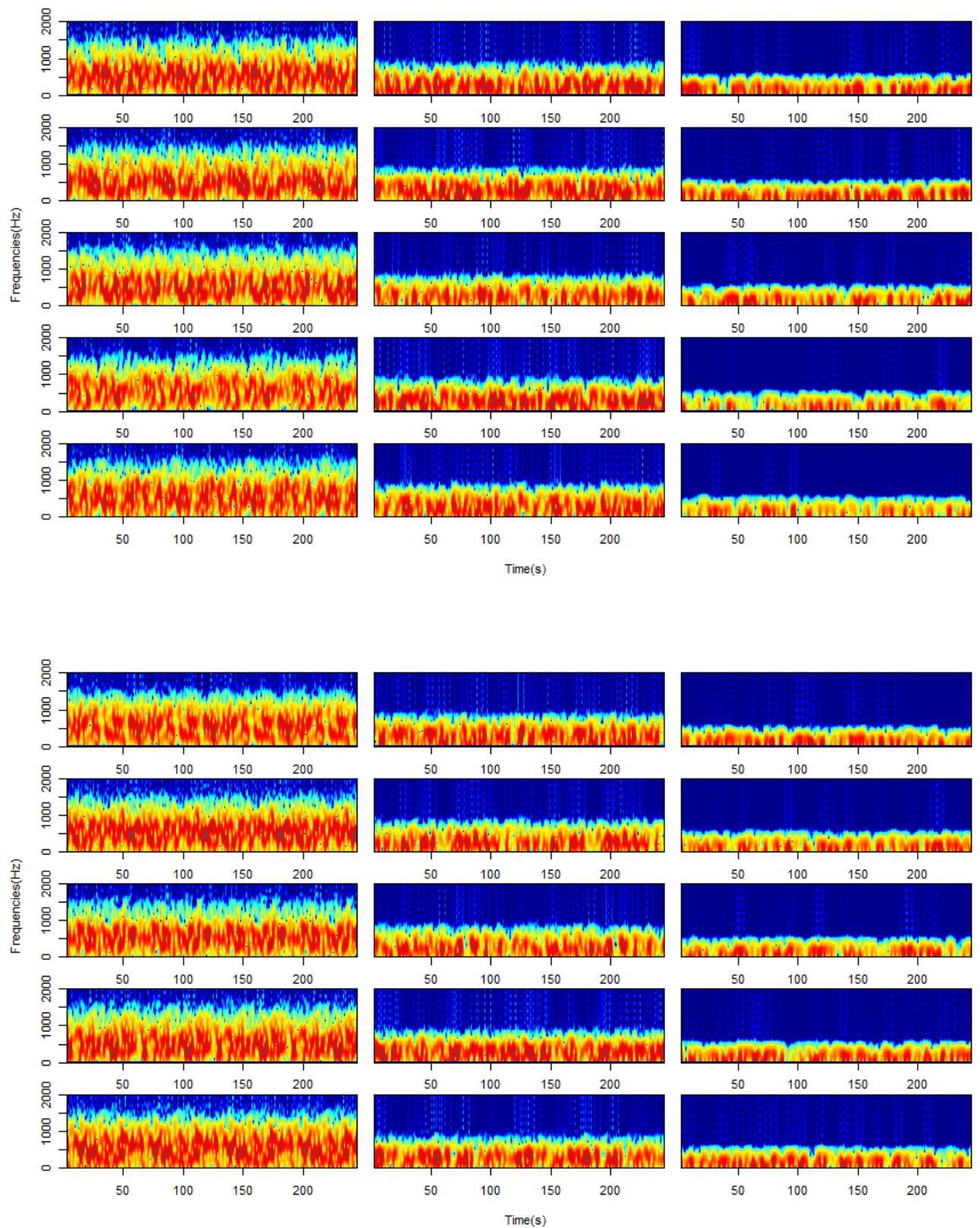


Figure 3:  $N = 1000, l = 0.05$ . Spectrograms of the  $y^{(m)}$  constructed with the above matrices for  $m = 1, \dots, 10$ .

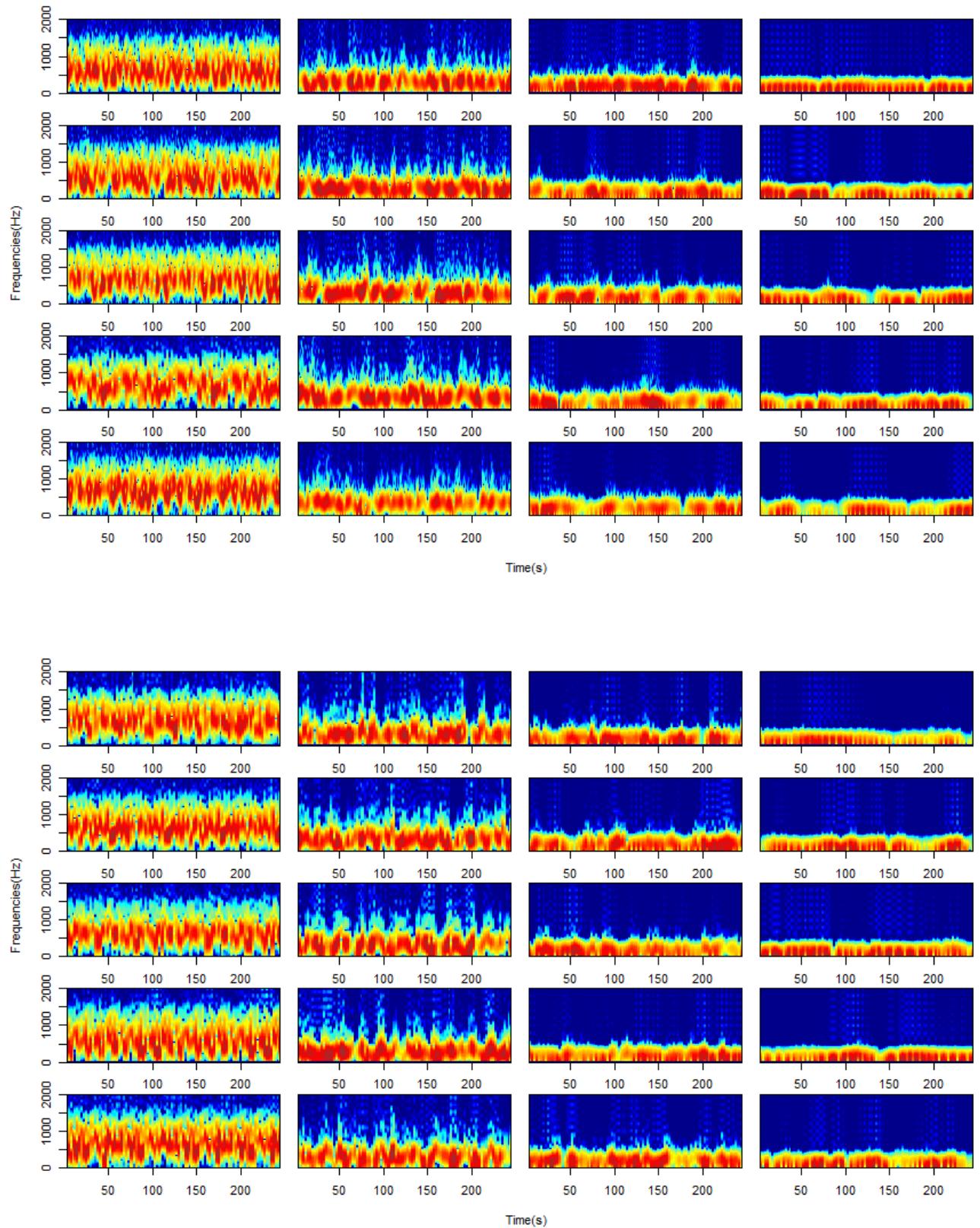


Figure 4:  $N = 1000, l = 0.05, p = 0.25, 0.5, 1$ . Spectrograms of the IMFs extracted by  $\hat{x}^{(m)}$  constructed with the above  $y_m$  for  $m = 1, \dots, 10$ .

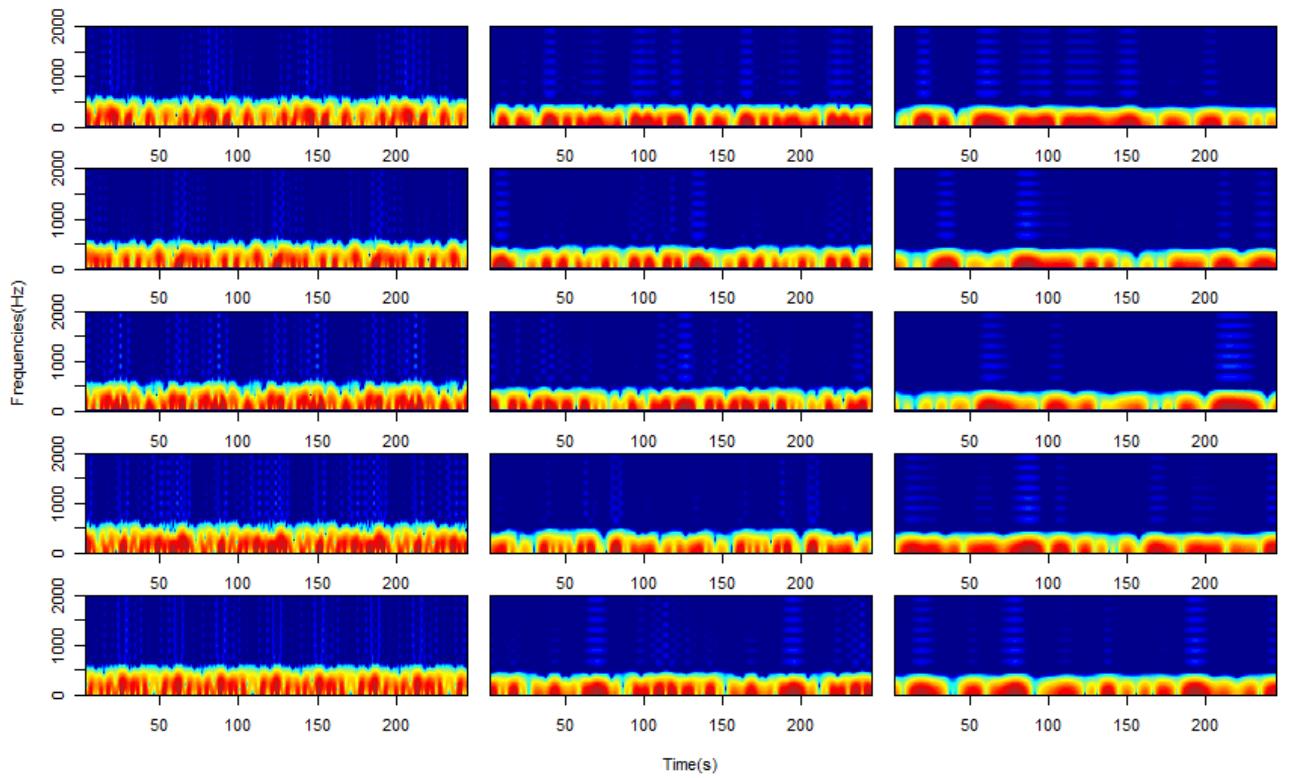
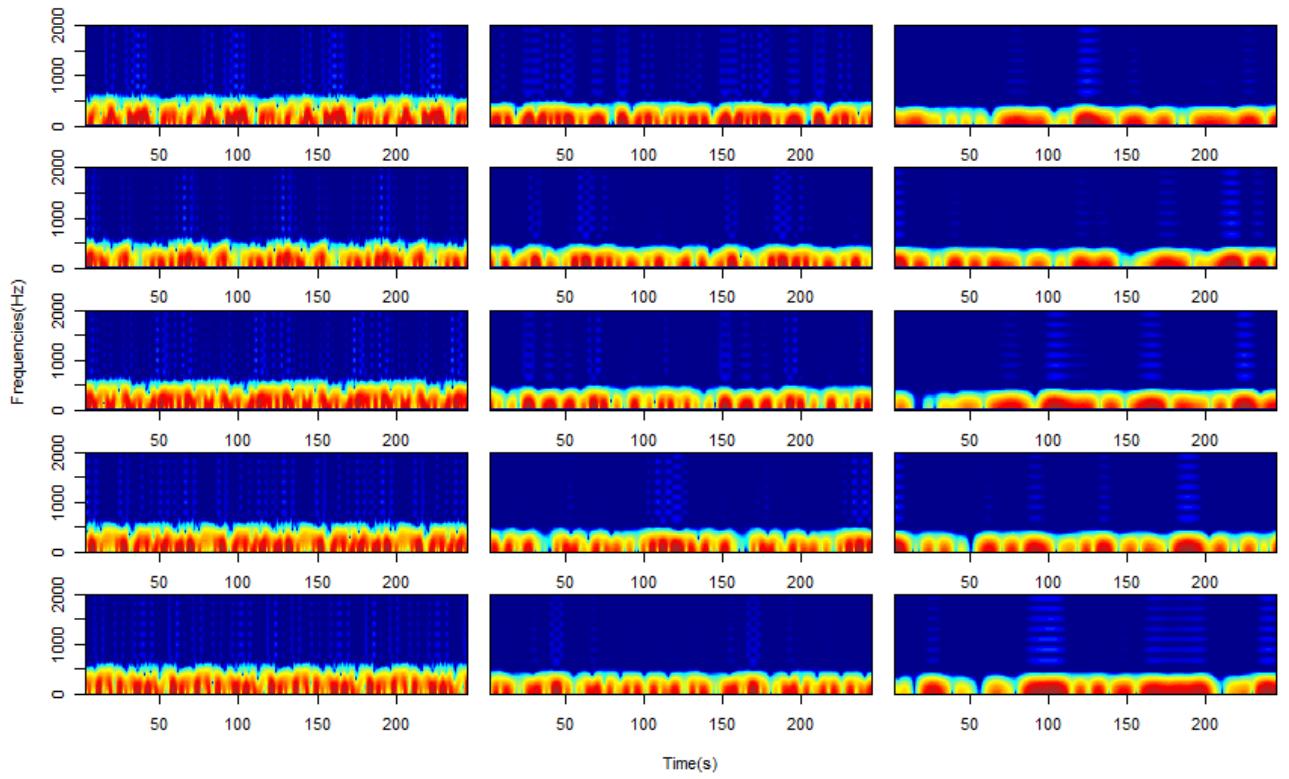


Figure 5:  $N = 1000, l = 0.2, p = 0.25, 0.5, 1$ . Spectrograms of the  $y^{(m)}$  components constructed with the above matrices for  $m = 1, \dots, 10$ .

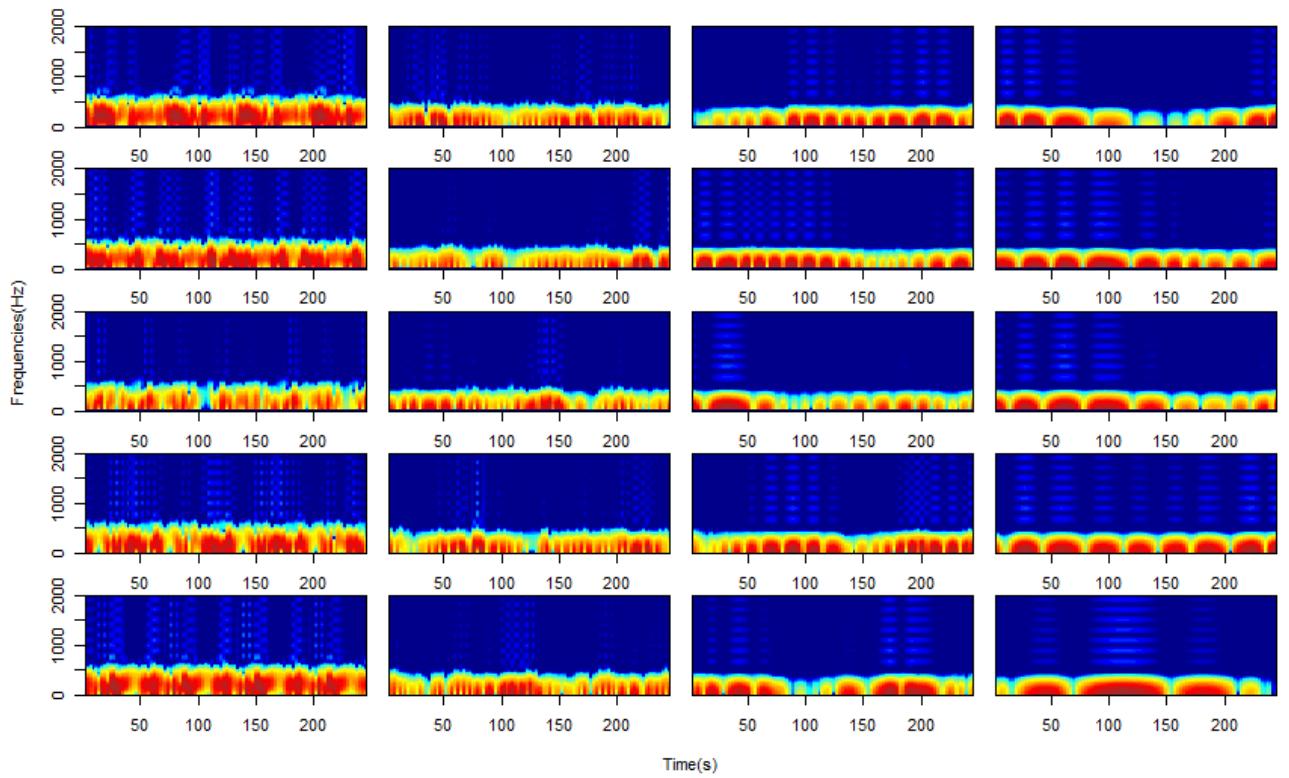
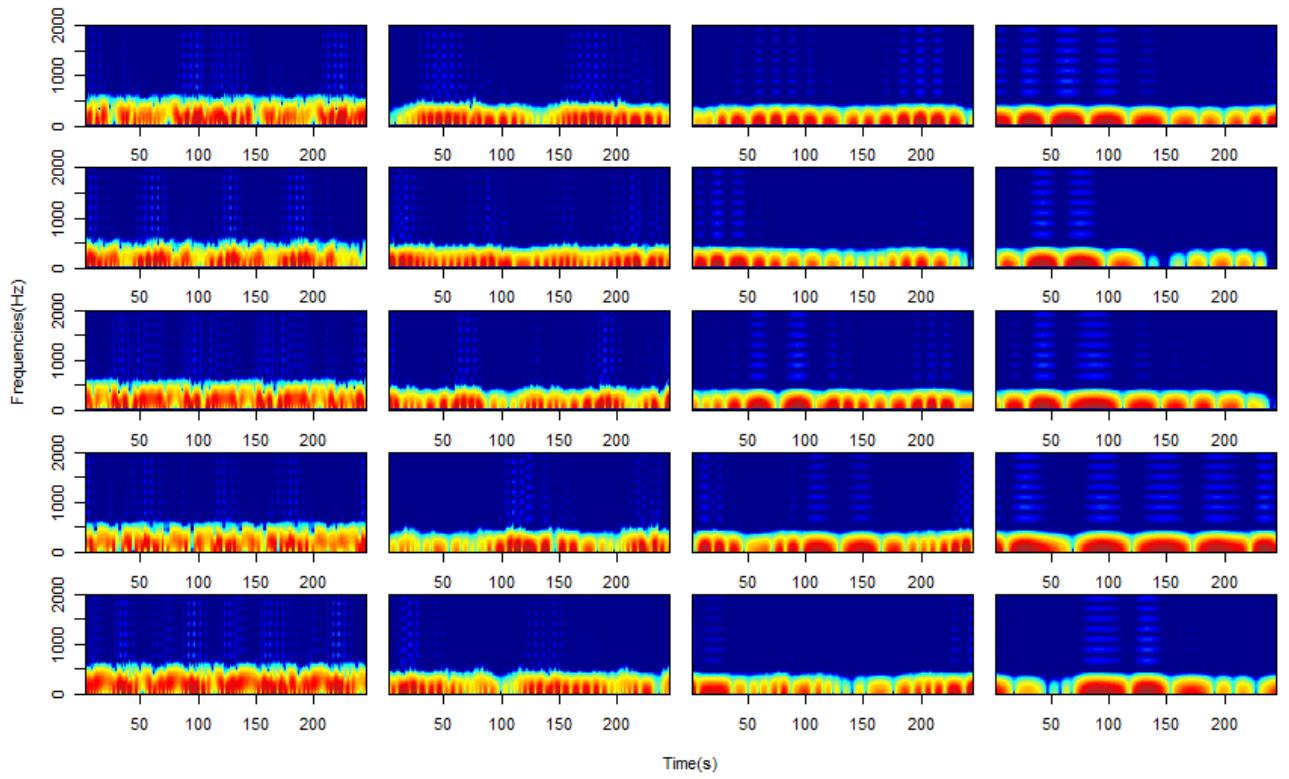


Figure 6:  $N = 1000, l = 0.2, p = 0.25, 0.5, 1$ . Spectrograms of the IMFs extracted by  $\hat{x}^{(m)}$  constructed with the above  $y^{(m)}$  for  $m = 1, \dots, 10$ .

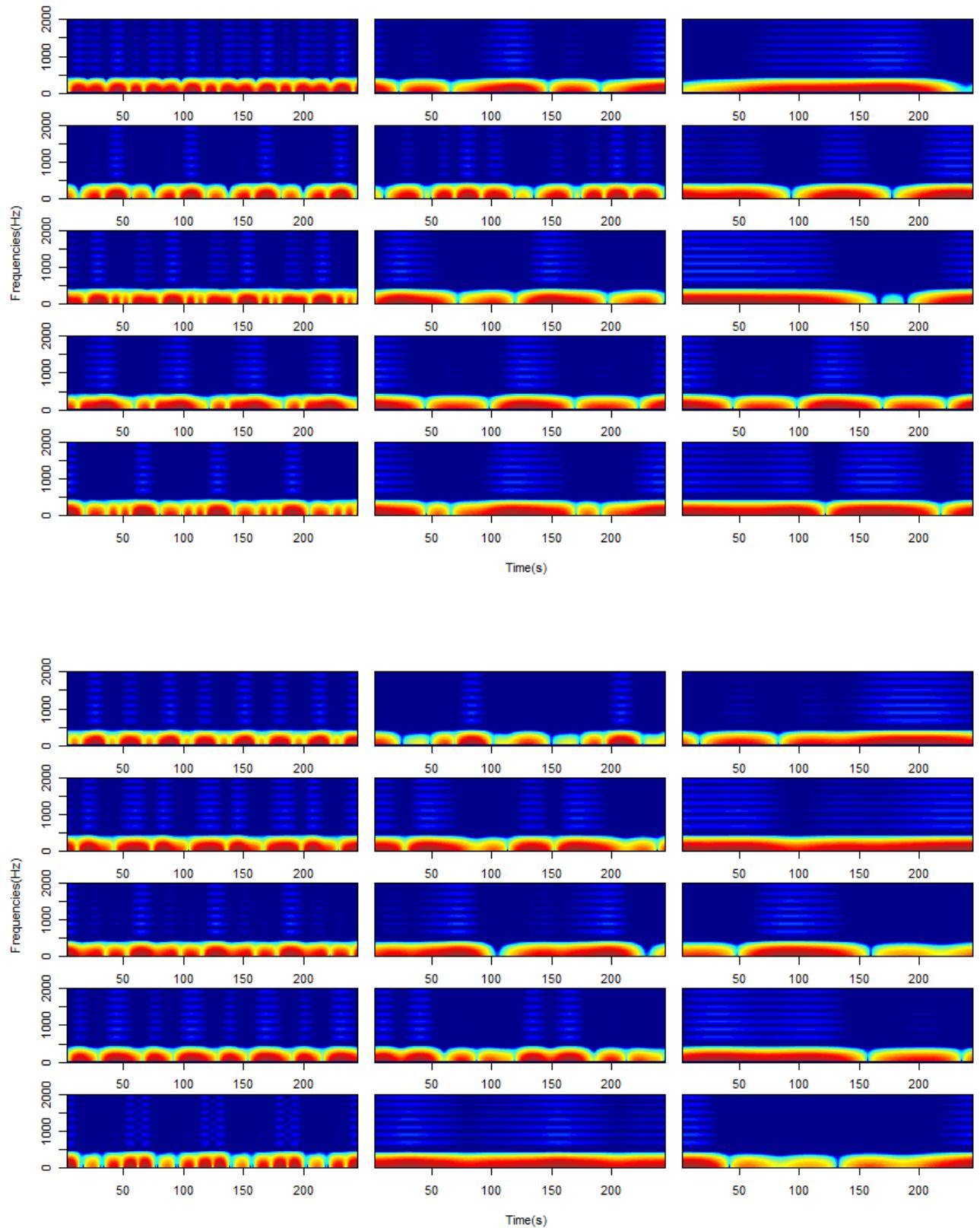


Figure 7:  $N = 1000, l = 1, p = 0.25, 0.5, 1$ . Spectrograms of the  $y^{(m)}$  components constructed with the above matrices for  $m = 1, \dots, 10$

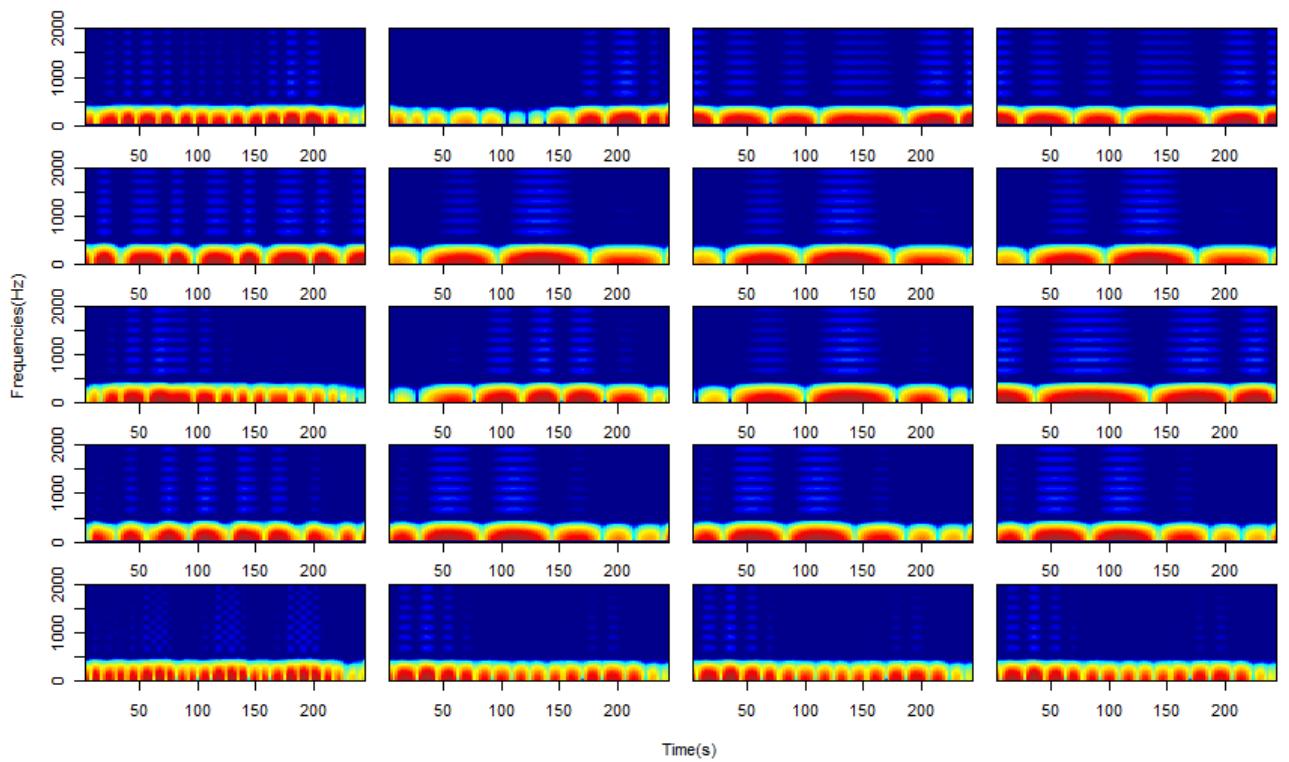
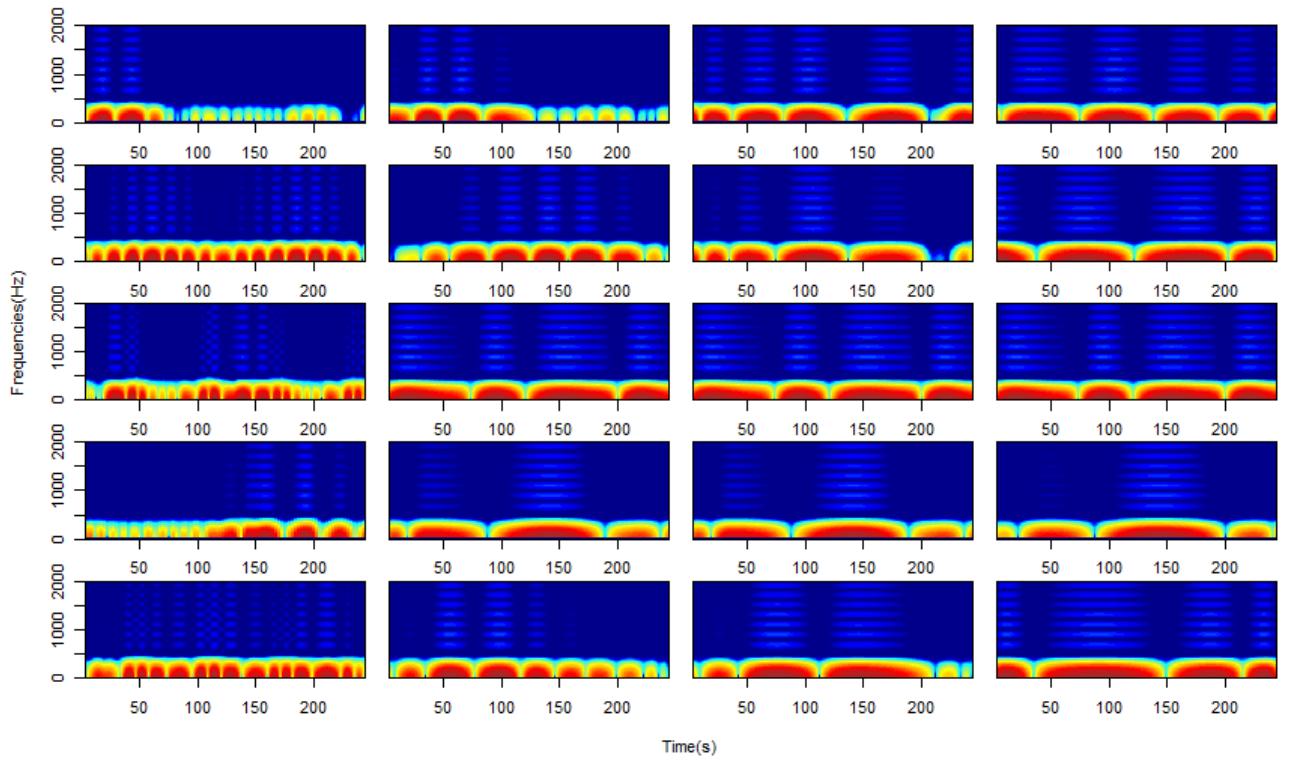


Figure 8:  $N = 1000, l = 1, p = 0.25, 0.5, 1, m = 1, \dots = 10$ . Spectrograms of the IMFs extracted by  $\hat{x}^{(m)}$  constructed with the above  $y^{(m)}$ .

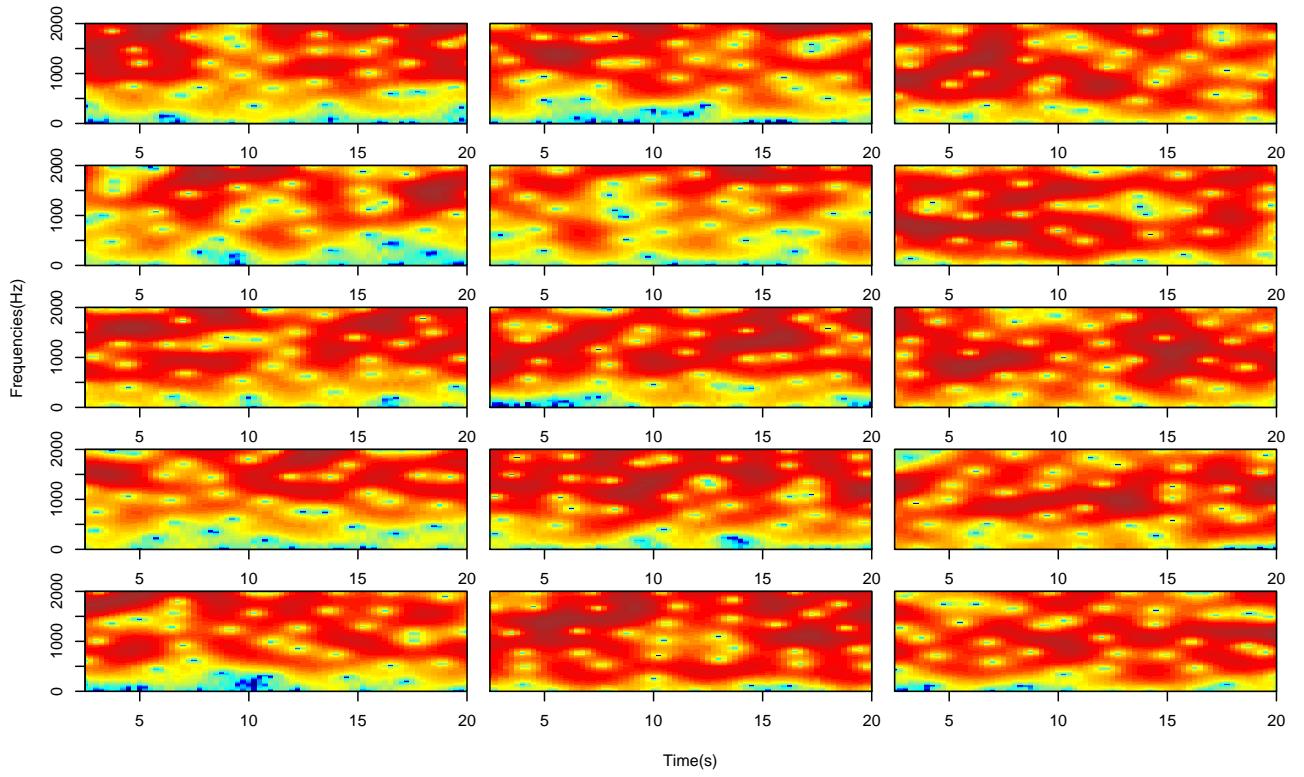
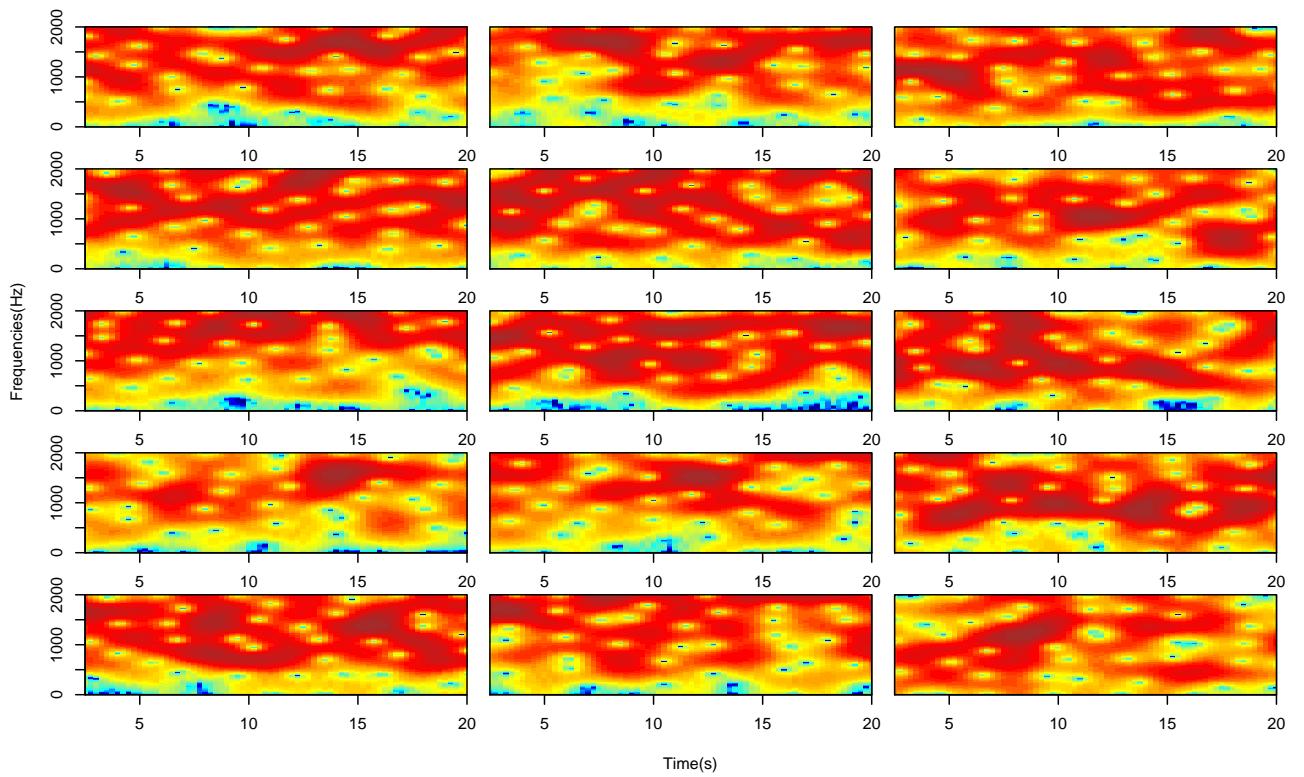


Figure 9:  $N = 100$ ,  $l = 0.05$ ,  $p = 0.25, 0.5, 1.00$ . Spectrograms of the  $y^{(m)}$  components.

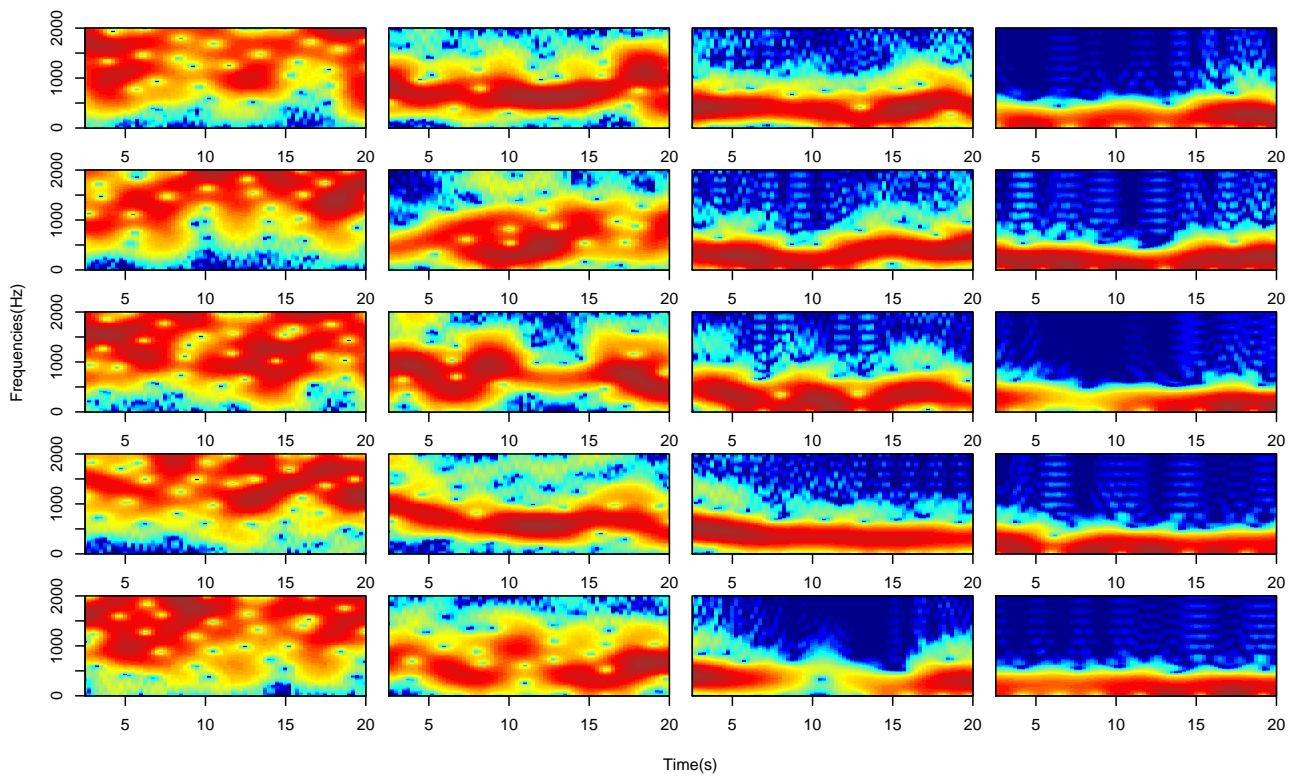
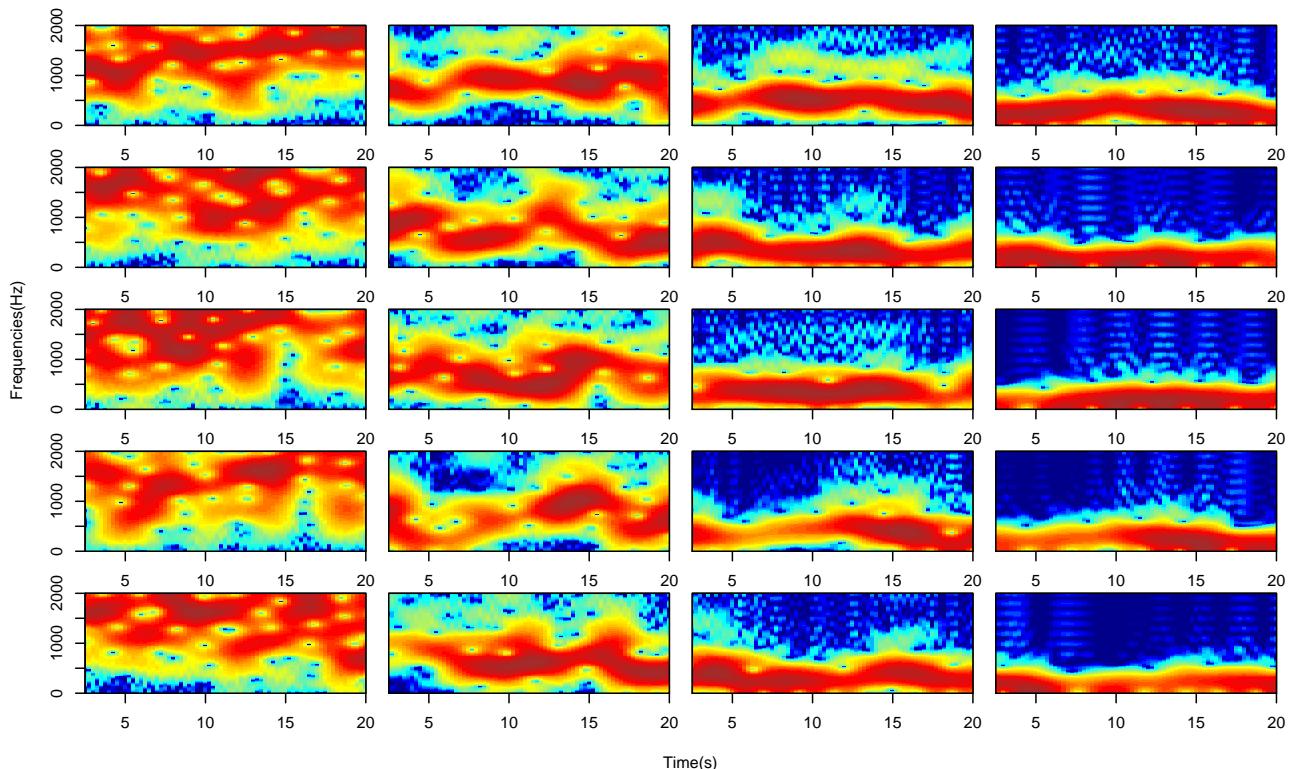


Figure 10:  $N = 100$ ,  $l = 0.05$ ,  $p = 0.25, 0.5, 1.00$ . Spectrograms of the IMFs extracted by  $\hat{x}^{(m)}$ .

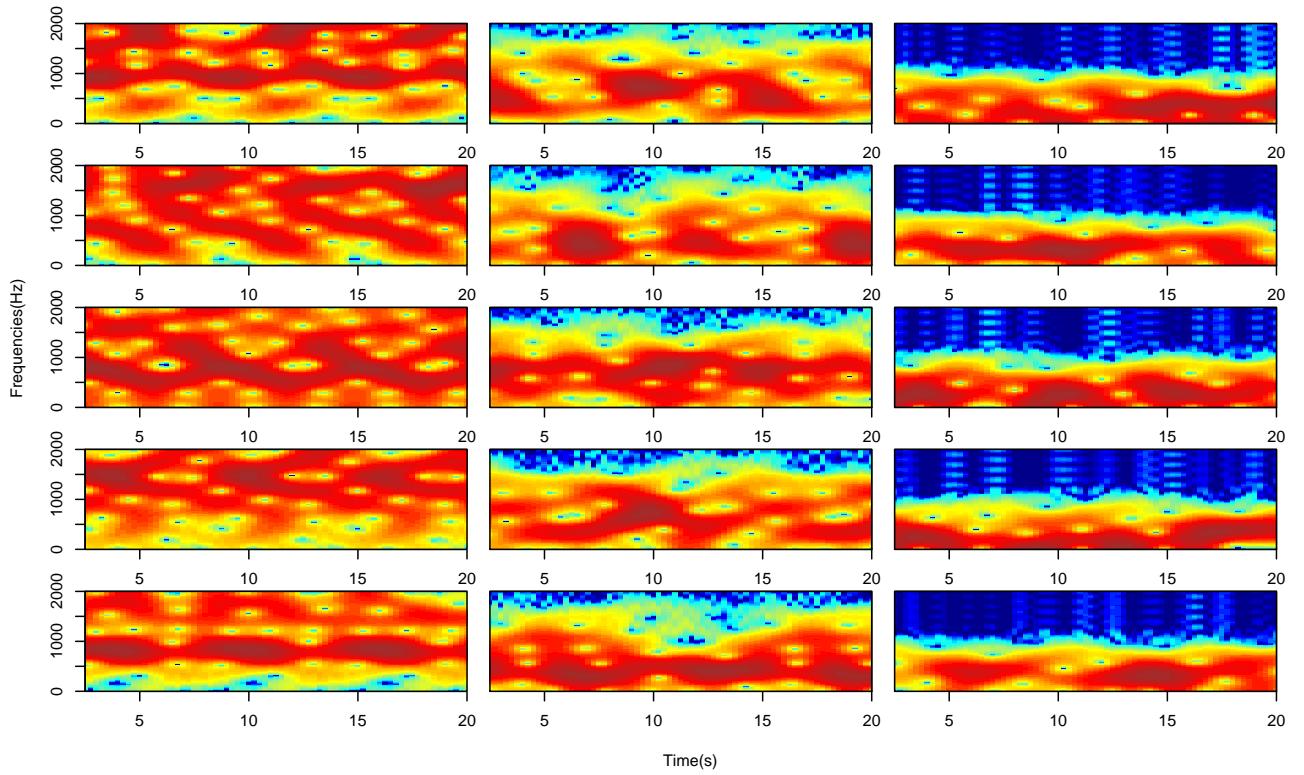
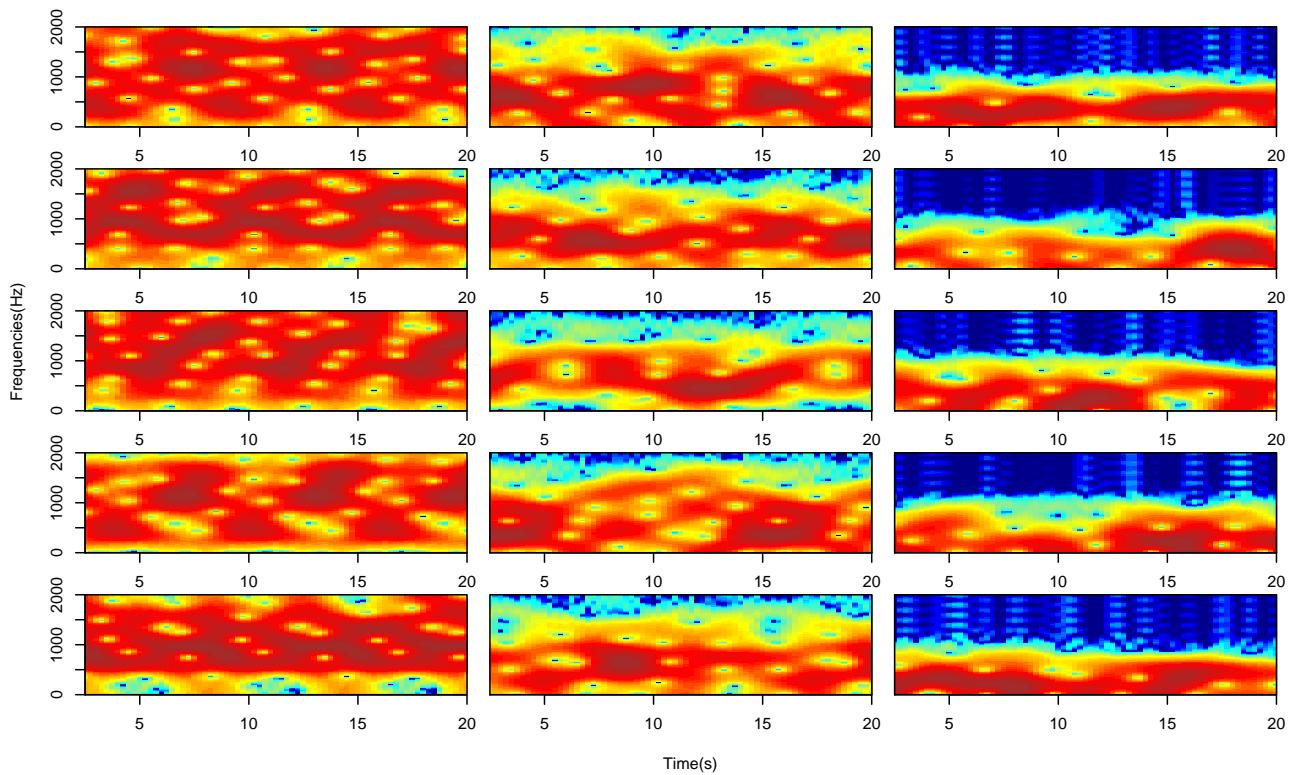


Figure 11:  $N = 100$ ,  $l = 0.2$ ,  $p = 0.25, 0.5, 1.00$ . Spectrograms of the  $y^{(m)}$  components.

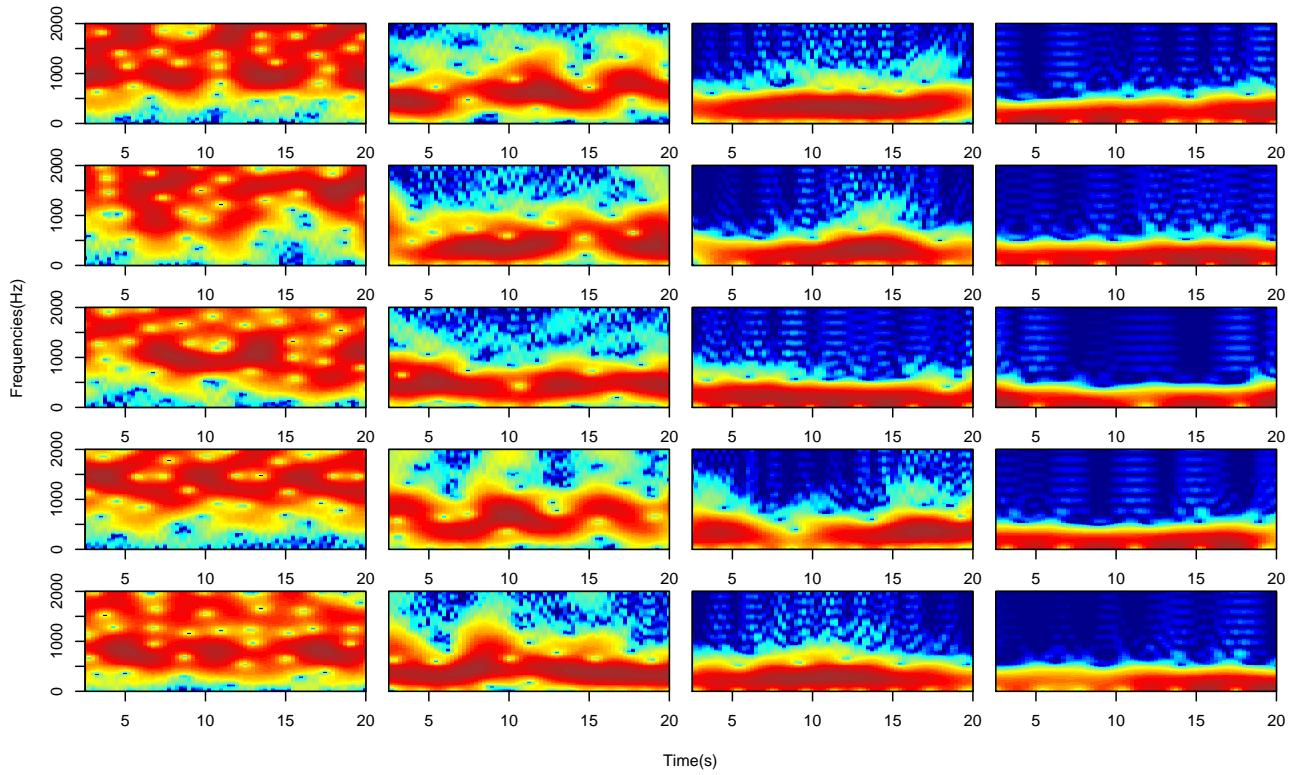
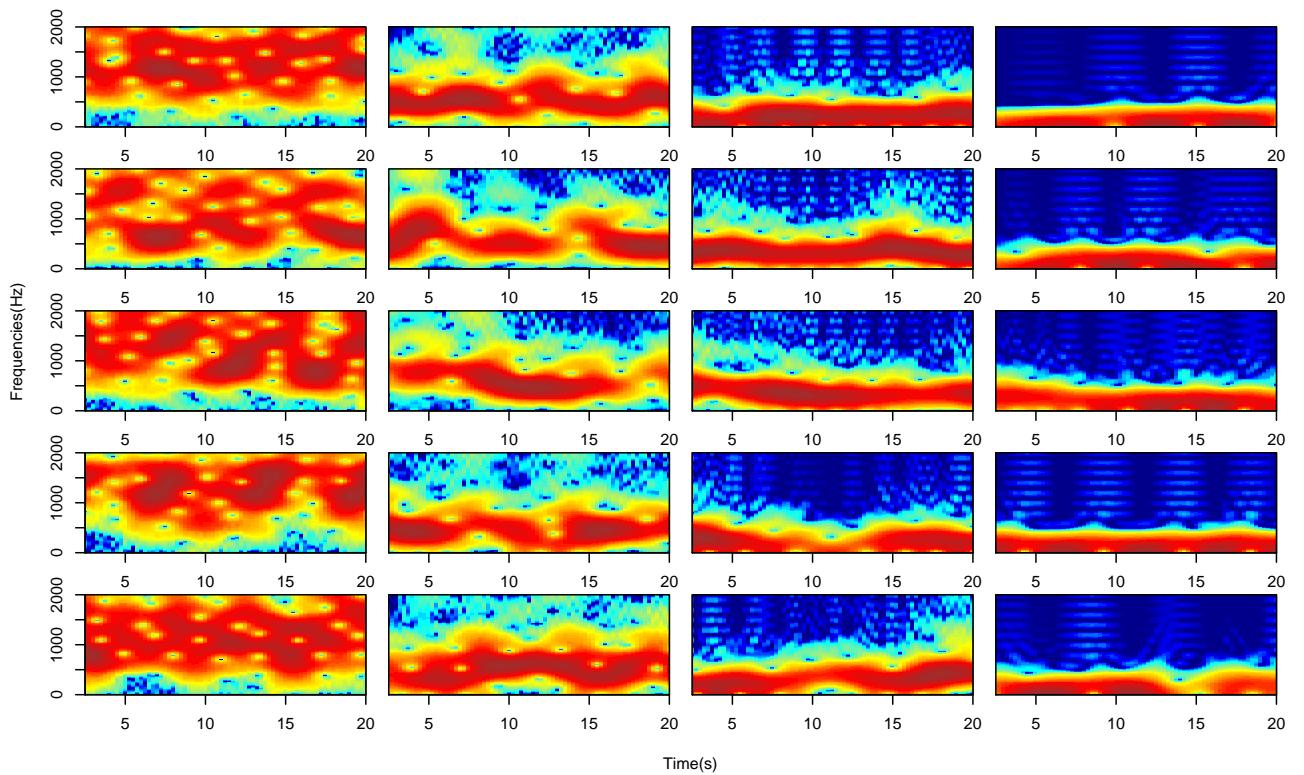


Figure 12:  $N = 100$ ,  $l = 0.2$ ,  $p = 0.25, 0.5, 1.00$ . Spectrograms of the IMFs extracted by  $\hat{x}^{(m)}$ .

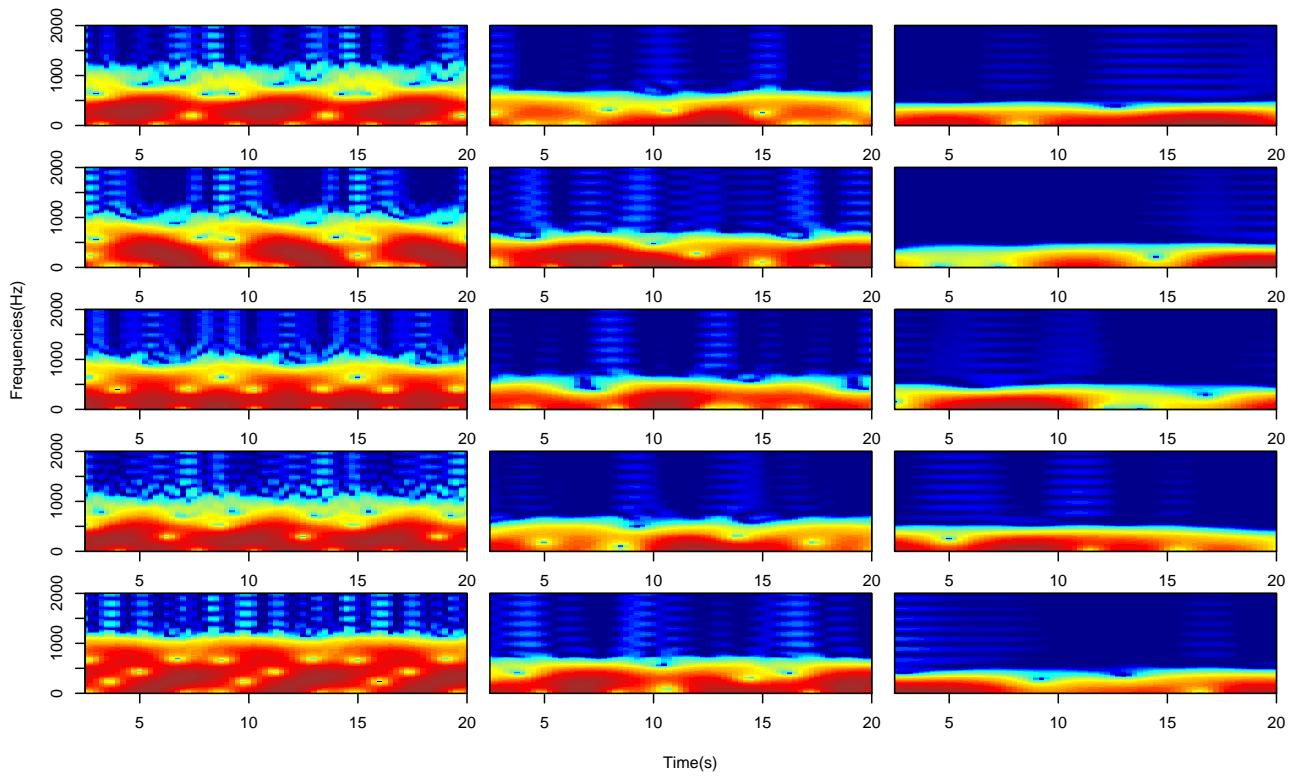
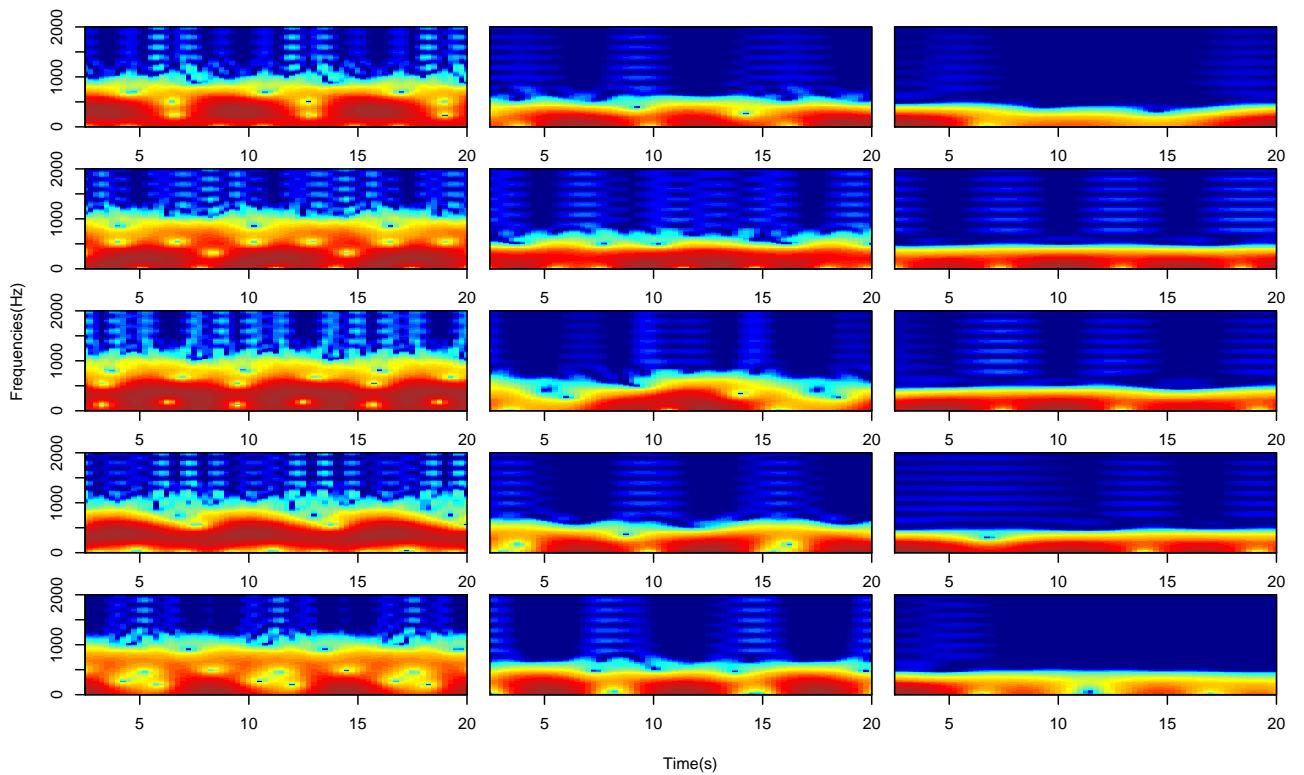


Figure 13:  $N = 100$ ,  $l = 1$ ,  $p = 0.25, 0.5, 1.00$ . Spectrograms of the  $y^{(m)}$  components.

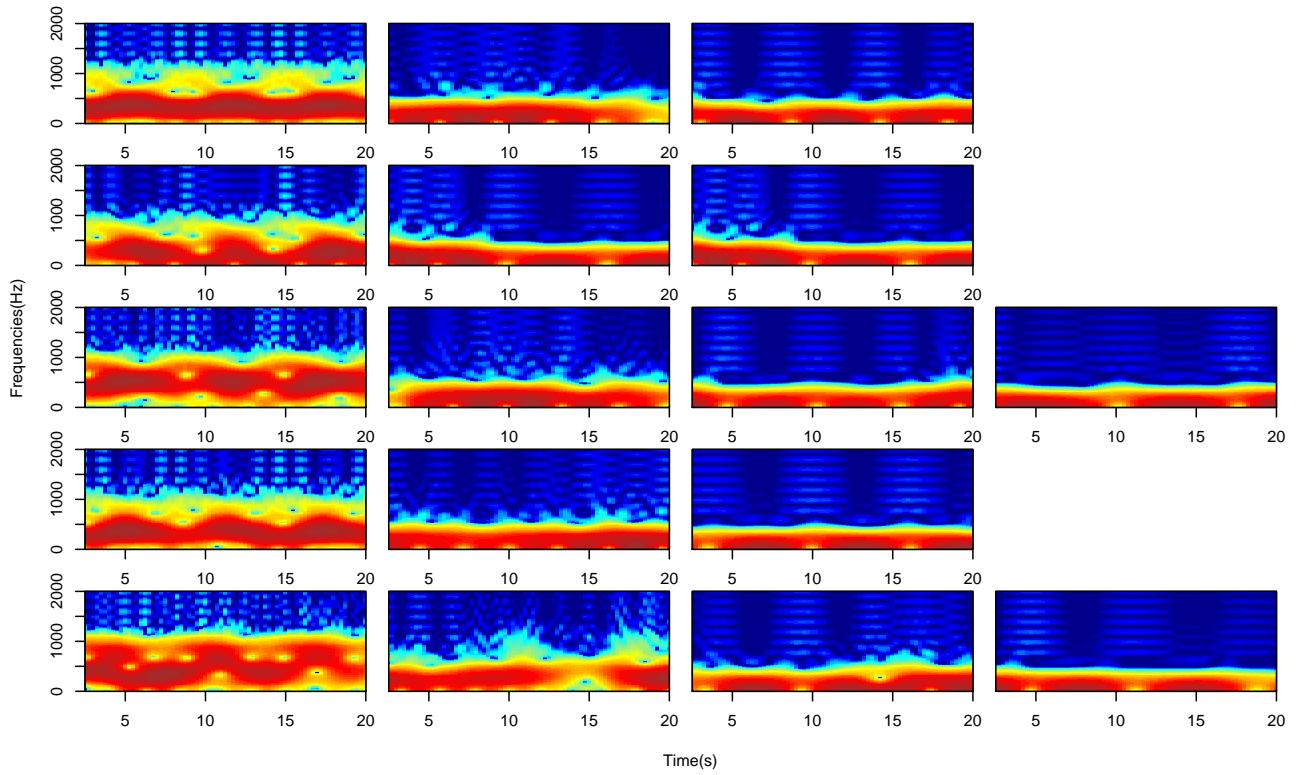
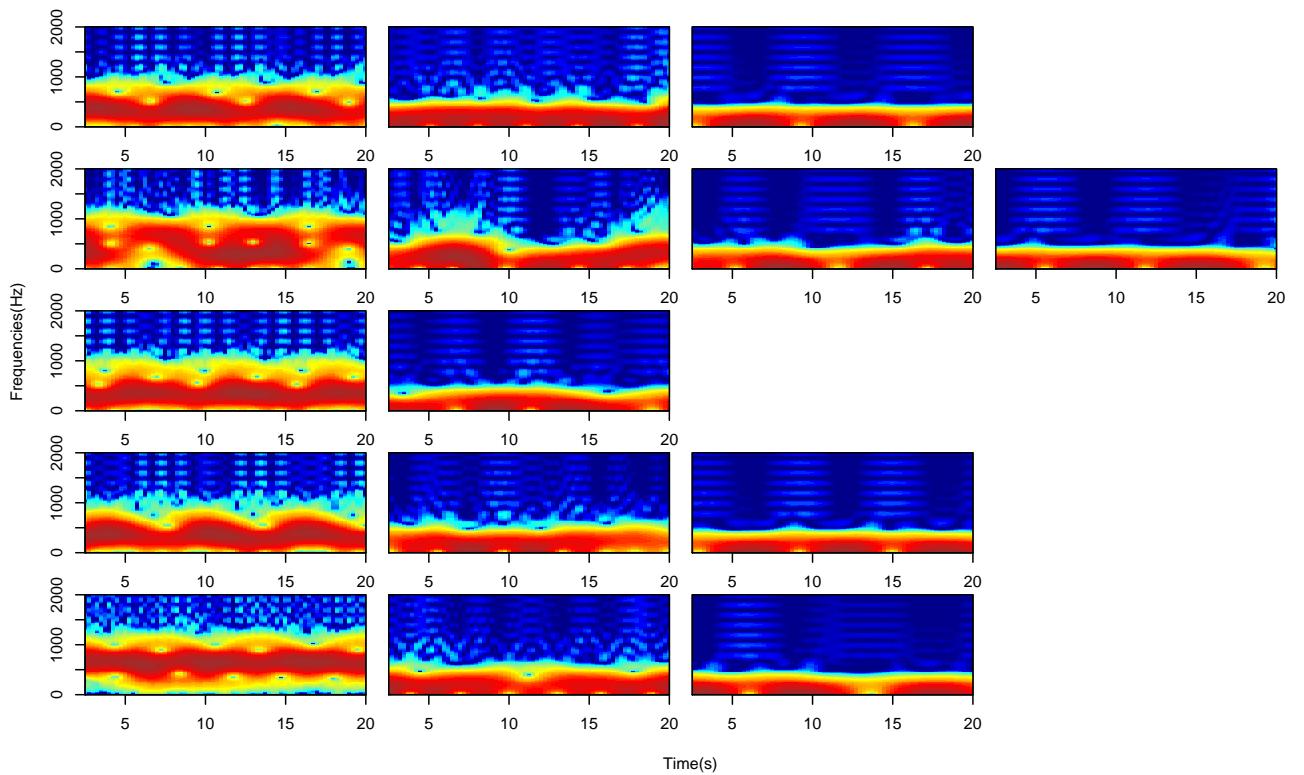


Figure 14:  $N = 100$ ,  $l = 1$ ,  $p = 0.25, 0.5, 1.00$ . Spectrograms of the IMFs extracted by  $\hat{x}^{(m)}$ .

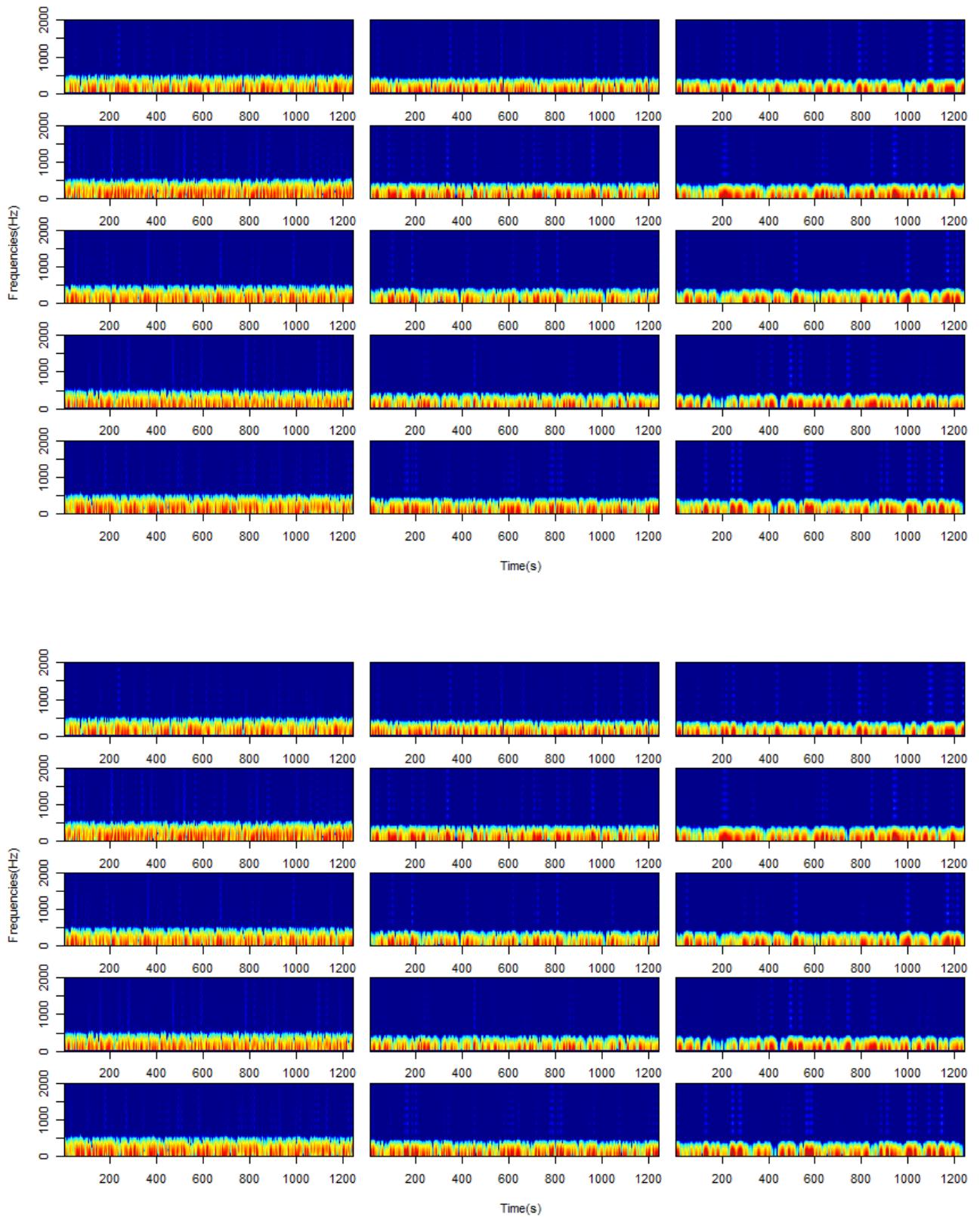


Figure 15:  $N = 5000$ ,  $l = 0.05$ ,  $p = 0.25, 0.5, 1.00$ . Spectrograms of the  $y^{(m)}$  components and IMFs extracted by  $\hat{x}^{(m)}$  for  $m = 1, 2, 3, 4, 5$ .

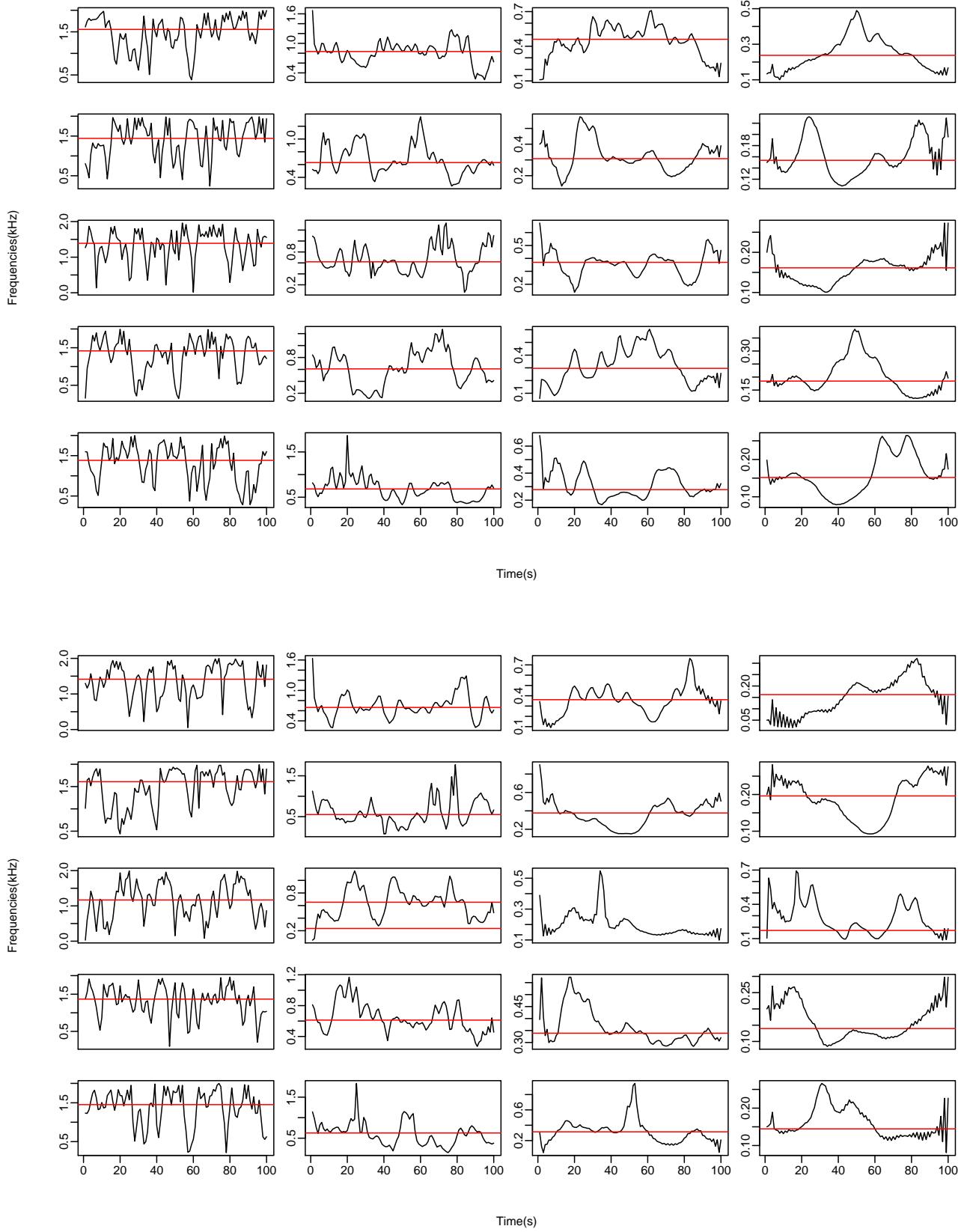


Figure 16:  $N = 100$ ,  $l = 0.05$ ,  $p = 0.25, 0.5, 1.00$ . Instantaneous Frequencies of the IMFs extracted by  $\hat{x}^{(m)}$ .

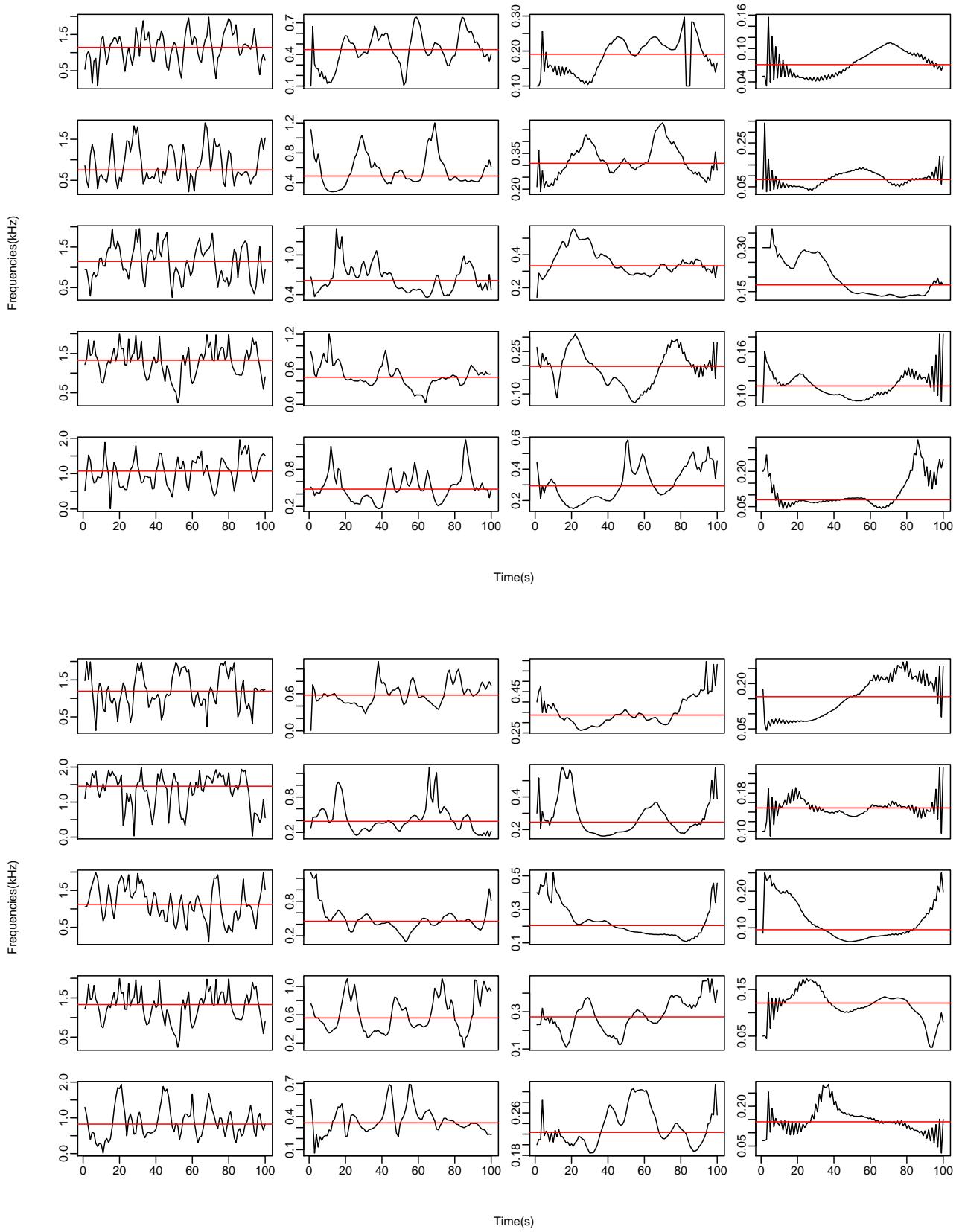


Figure 17:  $N = 100$ ,  $l = 0.2$ ,  $p = 0.25, 0.5, 1.00$ . Instantaneous Frequencies of the IMFs extracted by  $\hat{x}^{(m)}$ .

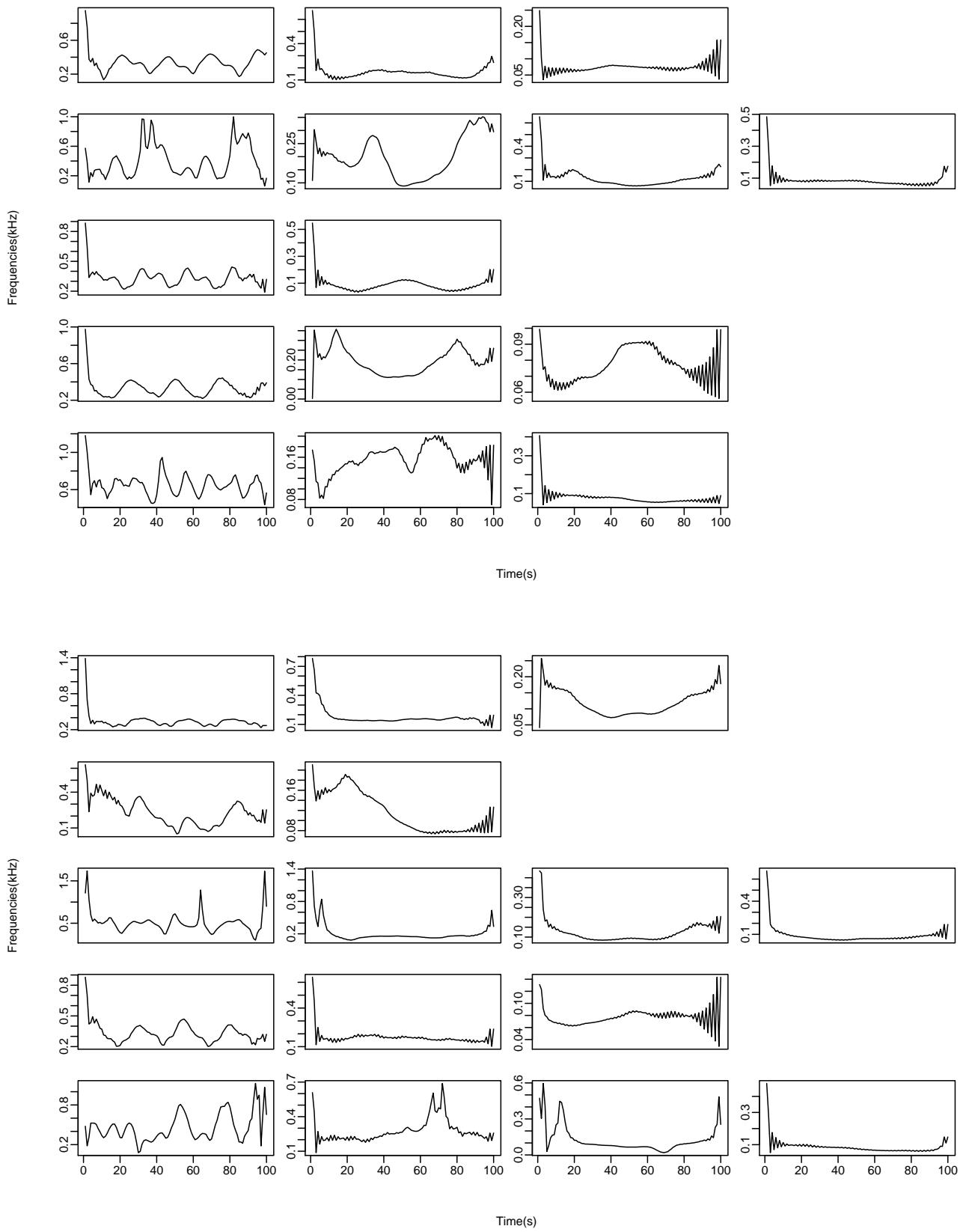


Figure 18:  $N = 100$ ,  $l = 1$ ,  $p = 0.25, 0.5, 1.00$ . Instantaneous Frequencies of the IMFs extracted by  $\hat{x}^{(m)}$ .

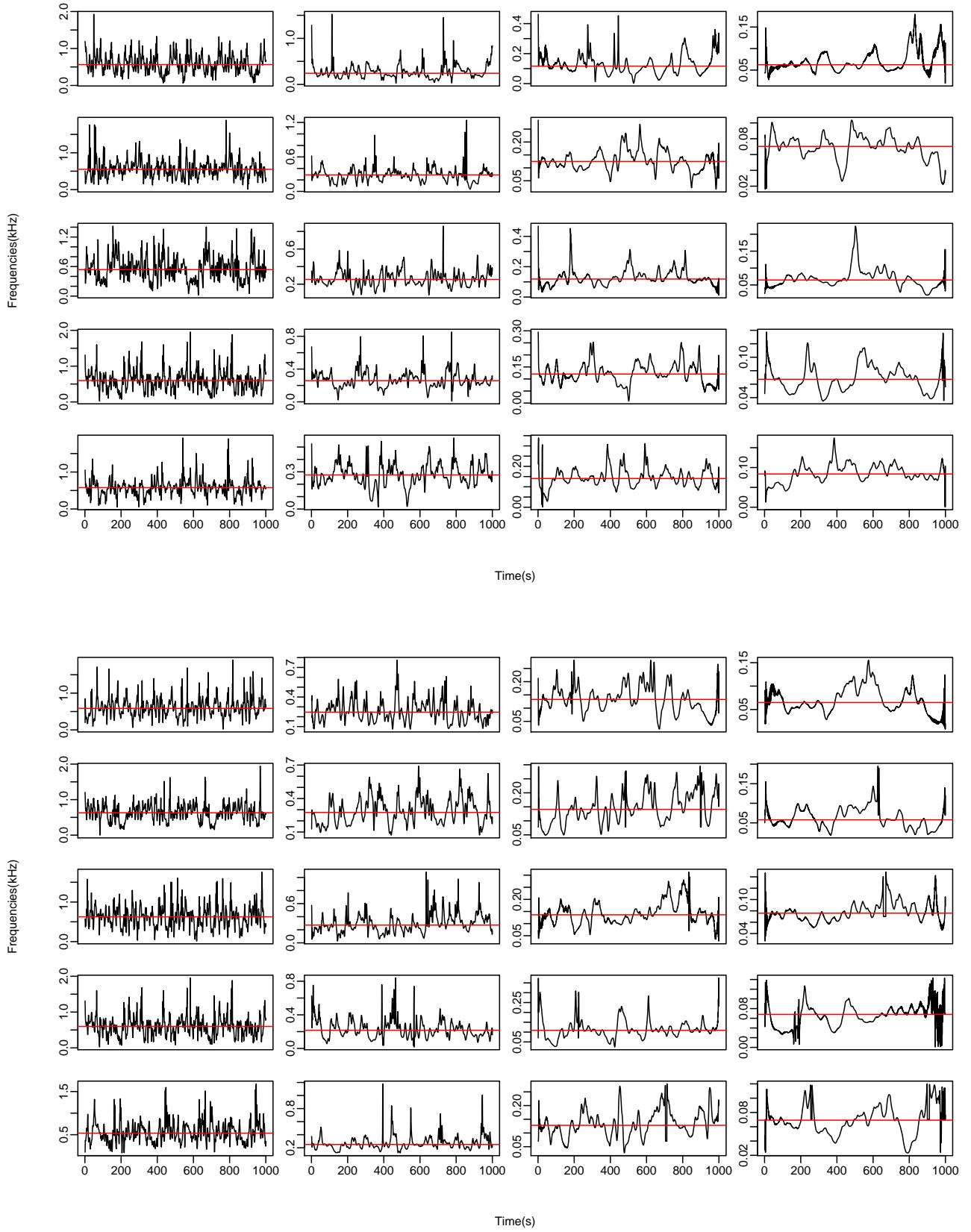


Figure 19:  $N = 1000$ ,  $l = 0.05$ ,  $p = 0.25, 0.5, 1.00$ . Instantaneous Frequencies of the IMFs extracted by  $\hat{x}^{(m)}$ .

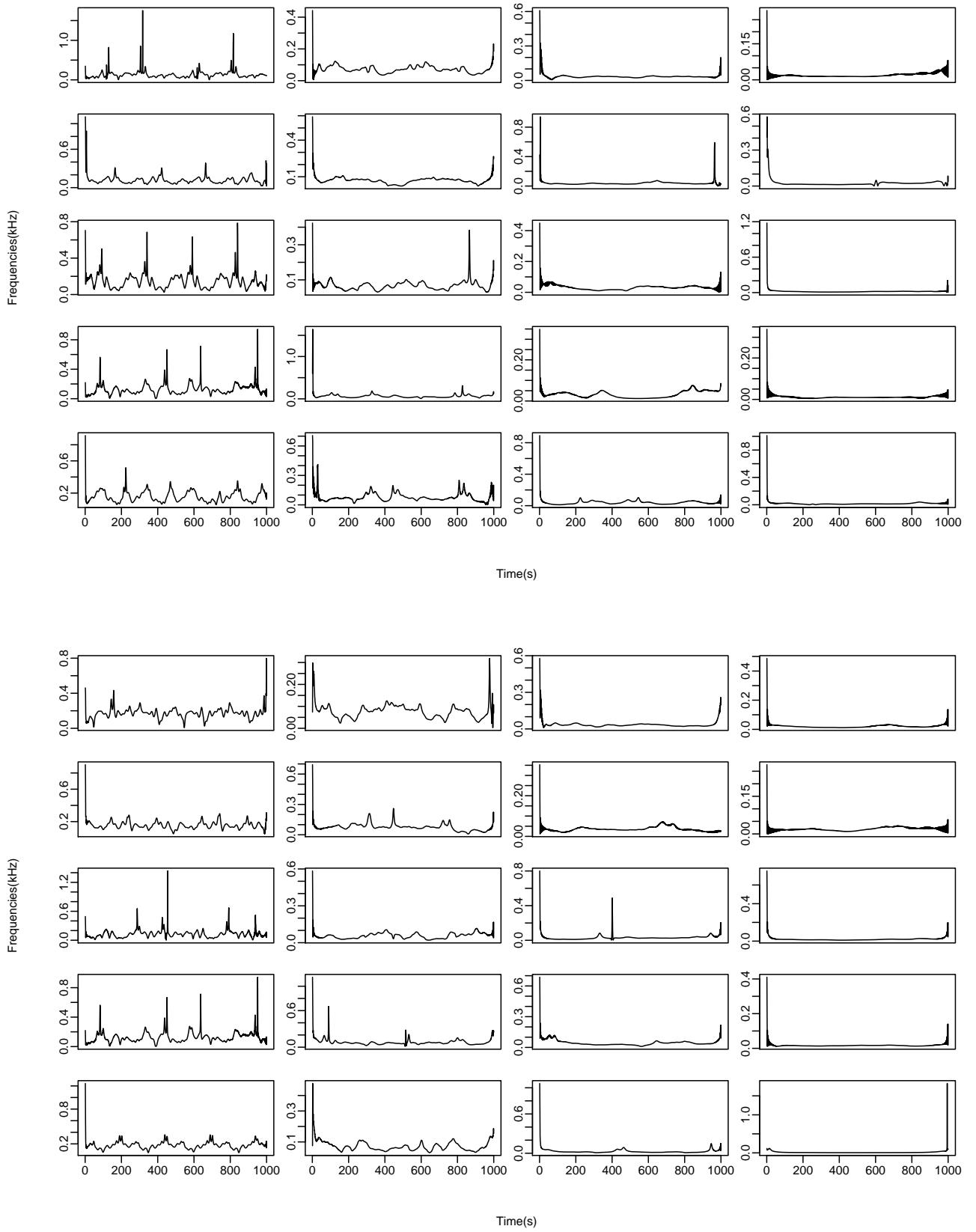


Figure 20:  $N = 1000$ ,  $l = 0.2$ ,  $p = 0.25, 0.5, 1.00$ . Instantaneous Frequencies of the IMFs extracted by  $\hat{x}^{(m)}$ .

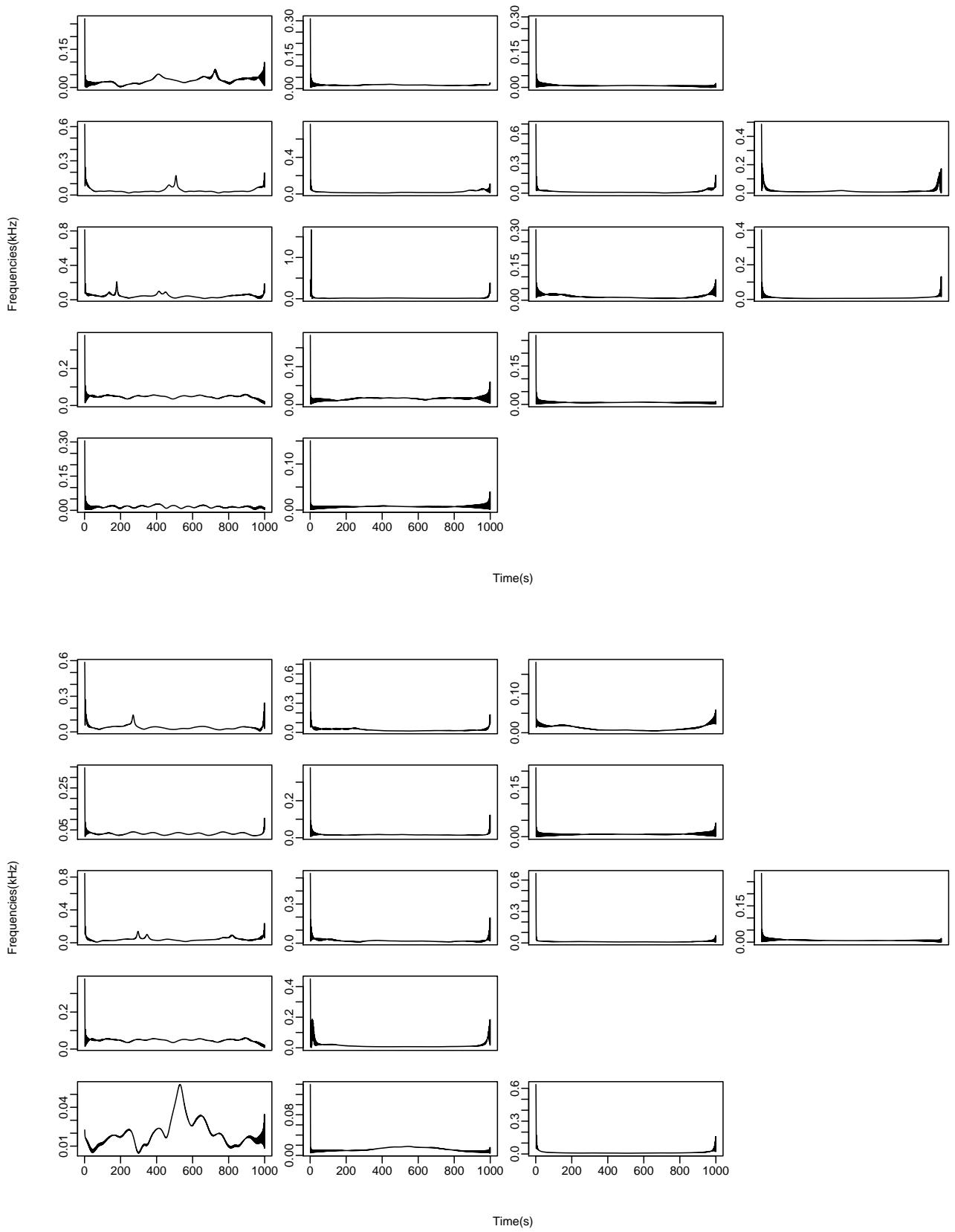


Figure 21:  $N = 1000$ ,  $l = 1$ ,  $p = 0.25, 0.5, 1.00$ . Instantaneous Frequencies of the IMFs extracted by  $\hat{x}^{(m)}$ .

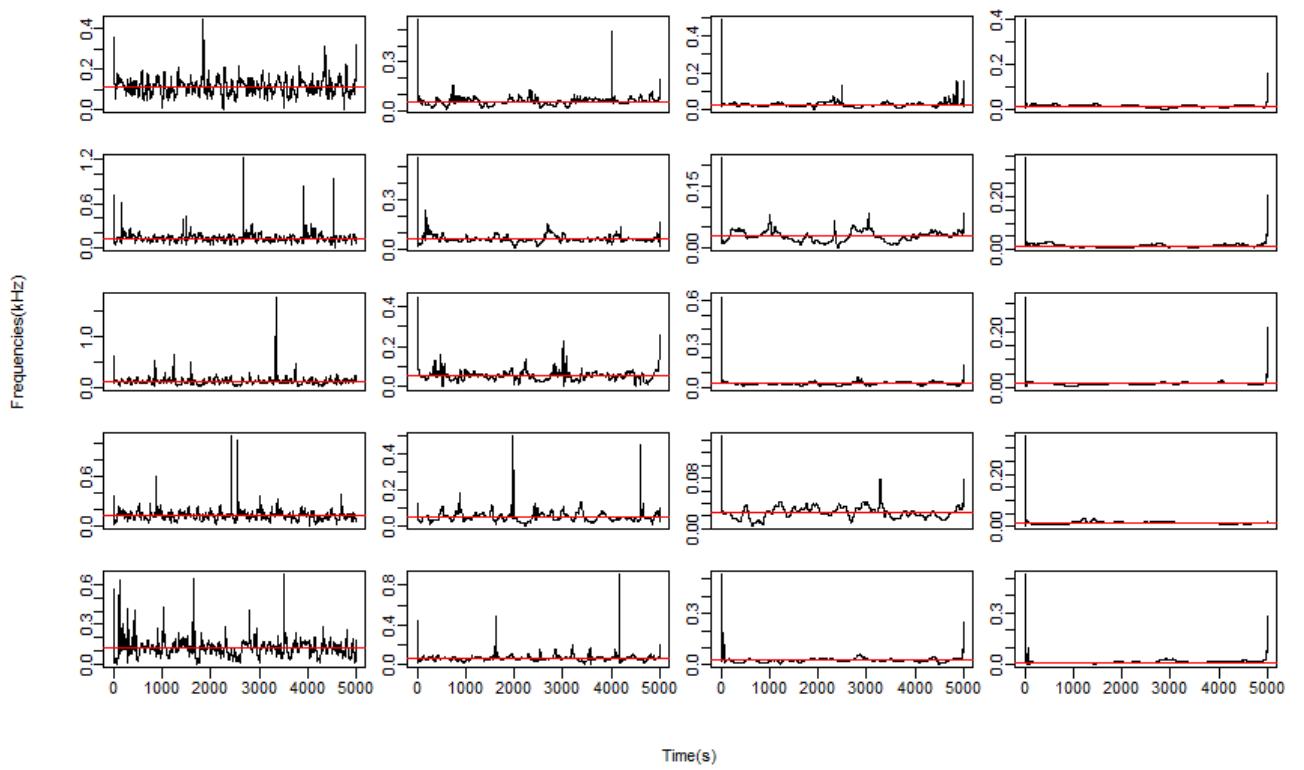


Figure 22:  $N = 5000$ ,  $l = 0.05$ ,  $p = 0.25, 0.5, 1.00$ . Instantaneous Frequencies of the IMFs extracted by  $\hat{x}^{(m)}$ ,  $m = 1, 2, 3, 4, 5$ .

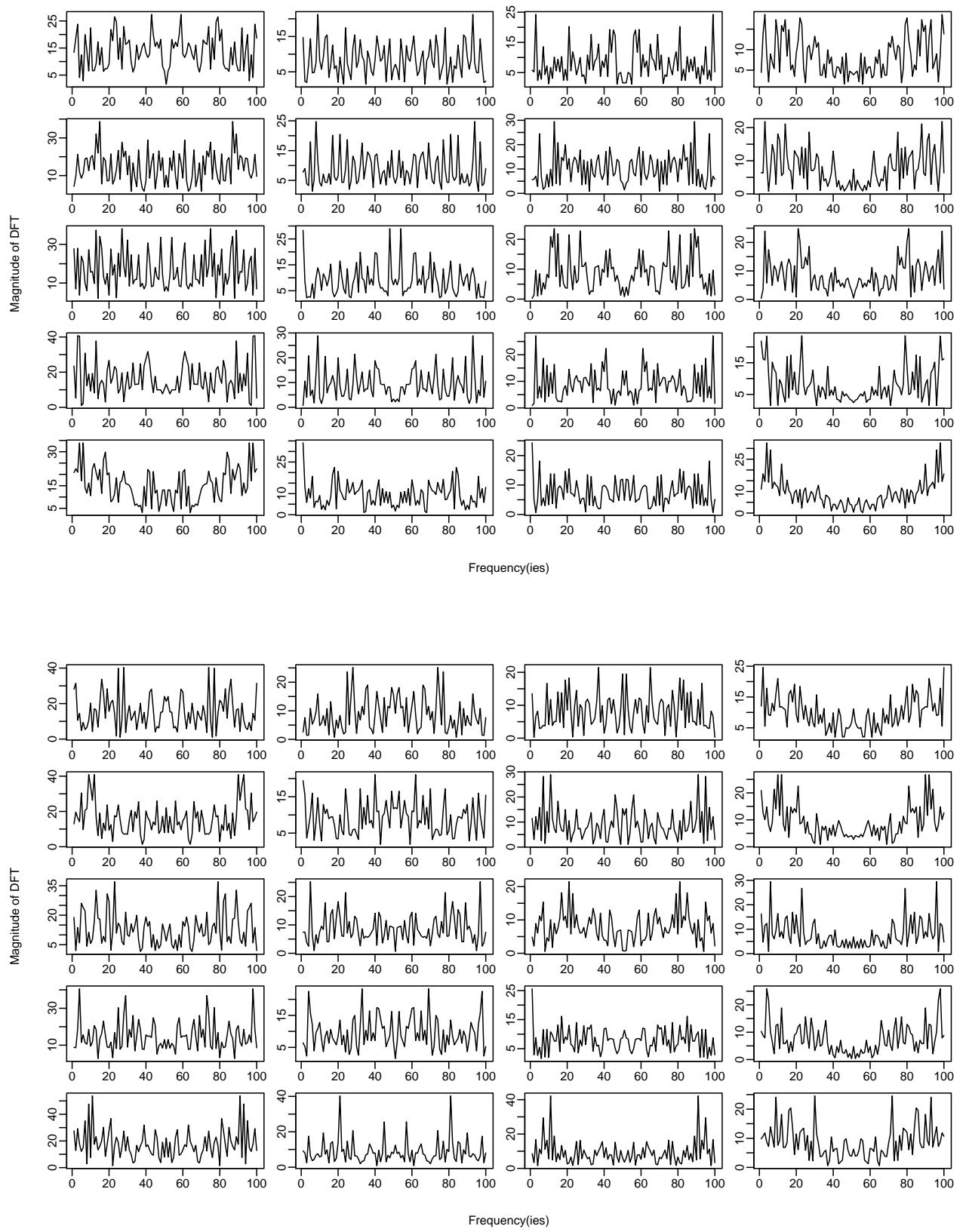


Figure 23:  $N = 100$ ,  $l = 0.05$ ,  $p = 0.25, 0.5, 1.00$ . Magnitudes of the DFT,  $m = 1, 2, \dots, 10$ .

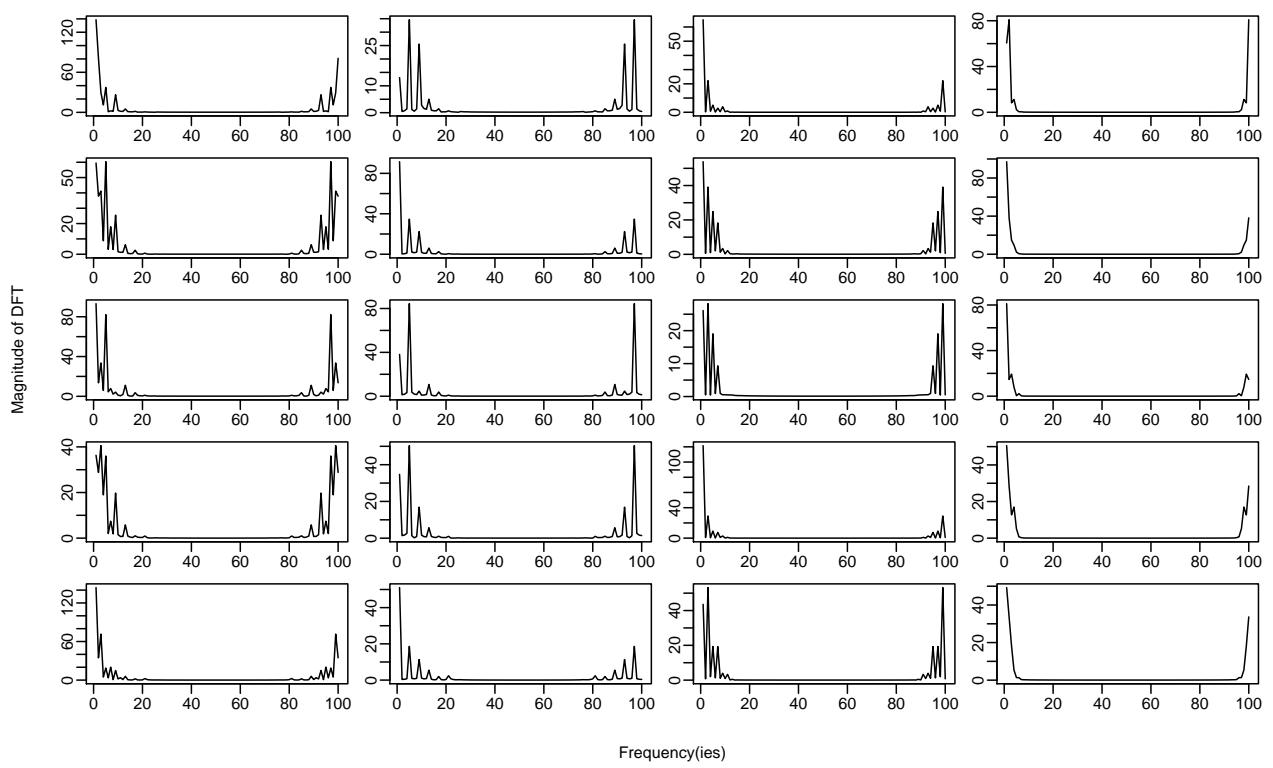
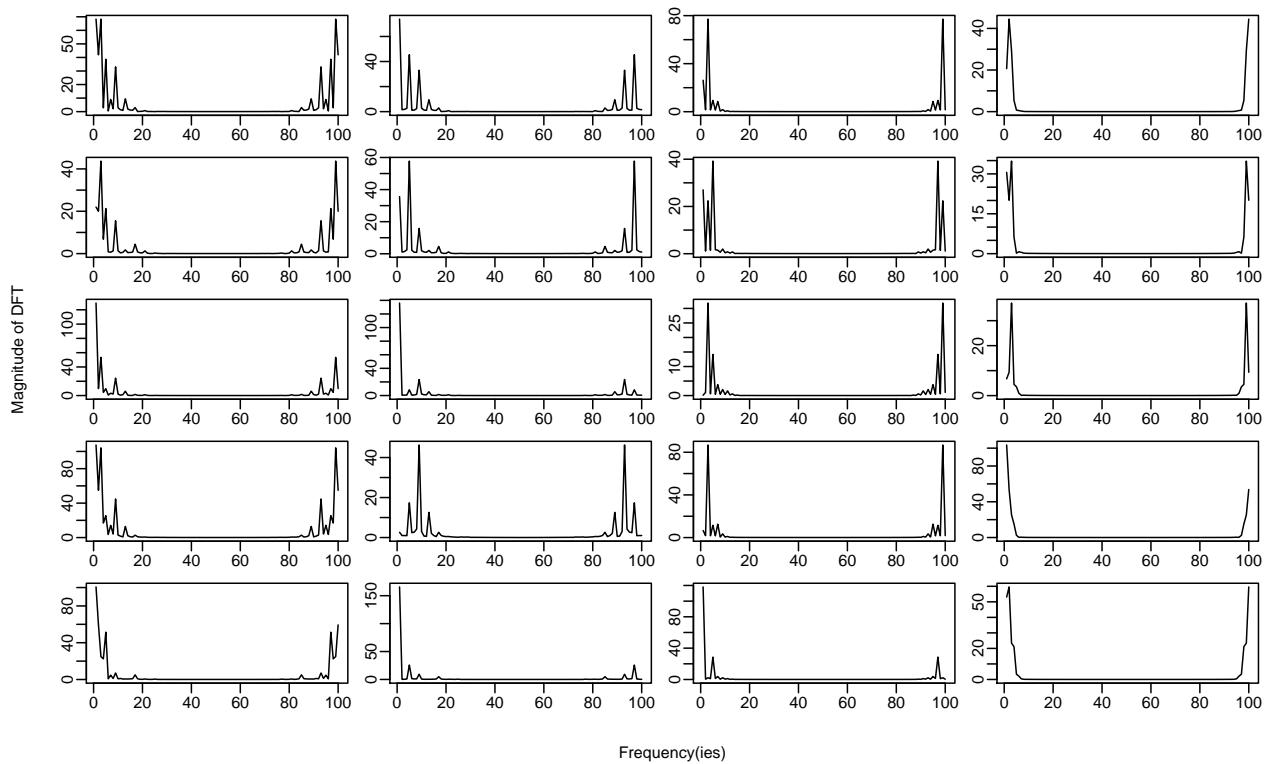


Figure 24:  $N = 100$ ,  $l = 1$ ,  $p = 0.25, 0.5, 1.00$ . Magnitudes of the DFT,  $m = 1, 2, \dots, 10$ .

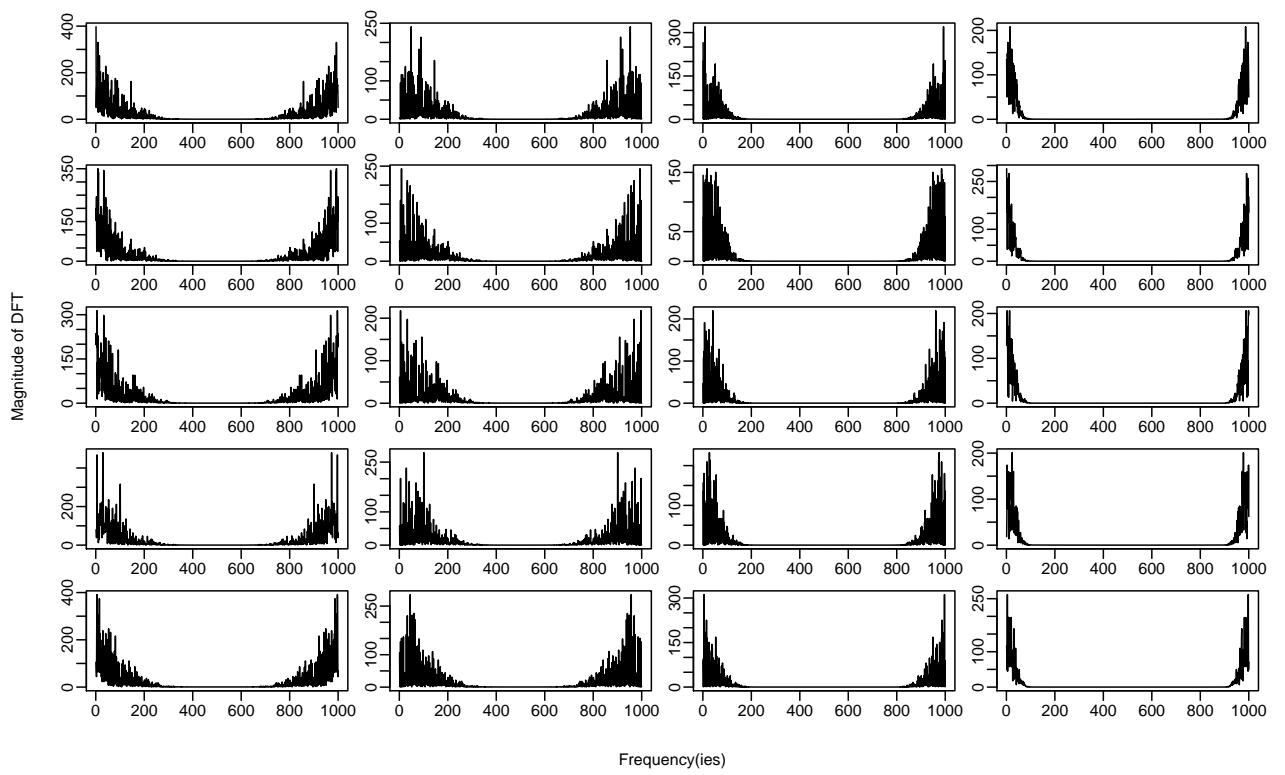
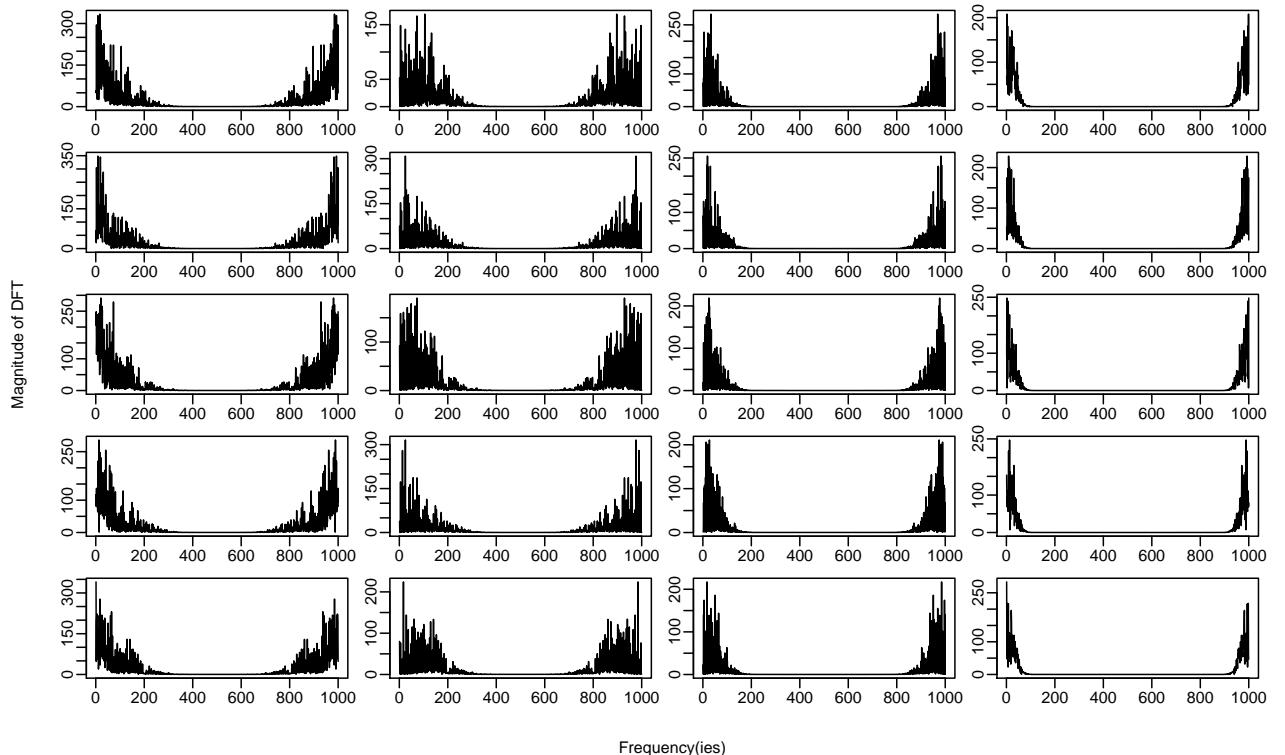


Figure 25:  $N = 1000$ ,  $l = 0.05$ ,  $p = 0.25, 0.5, 1.00$ . Magnitudes of the DFT,  $m = 1, 2, \dots, 10$ .

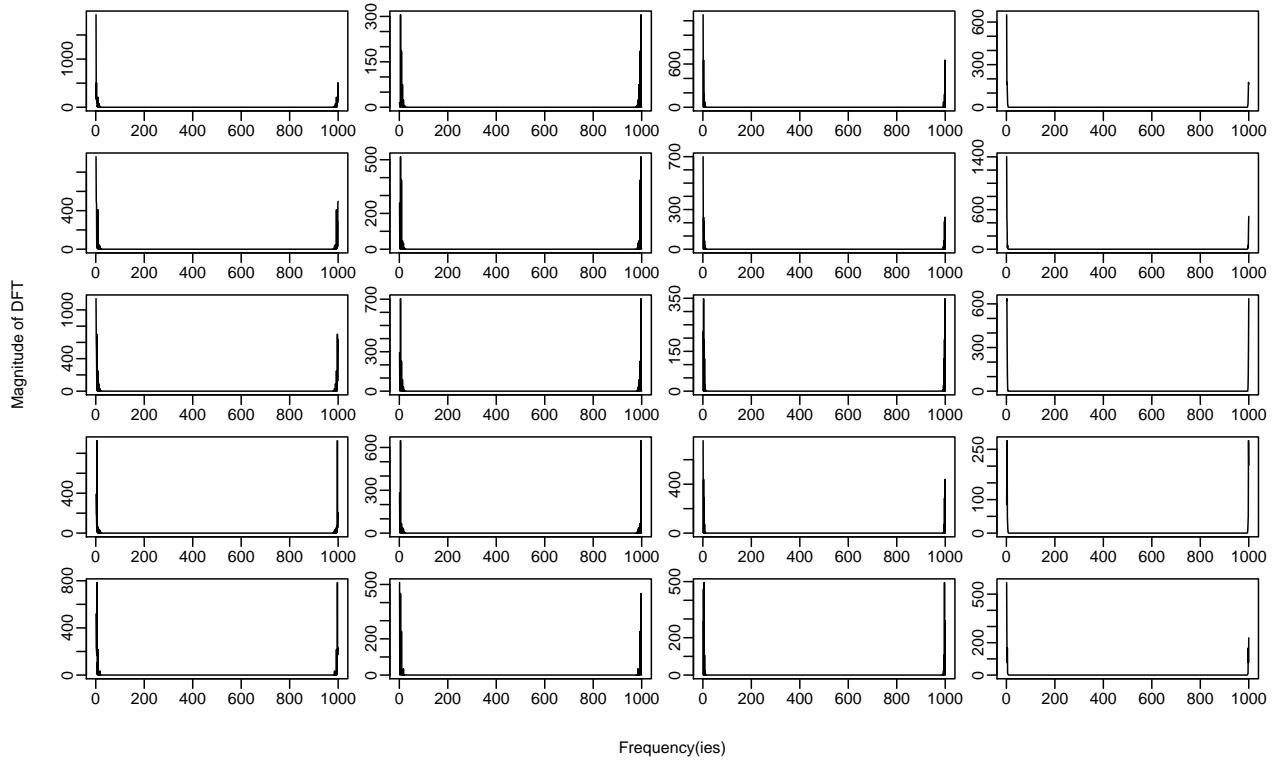
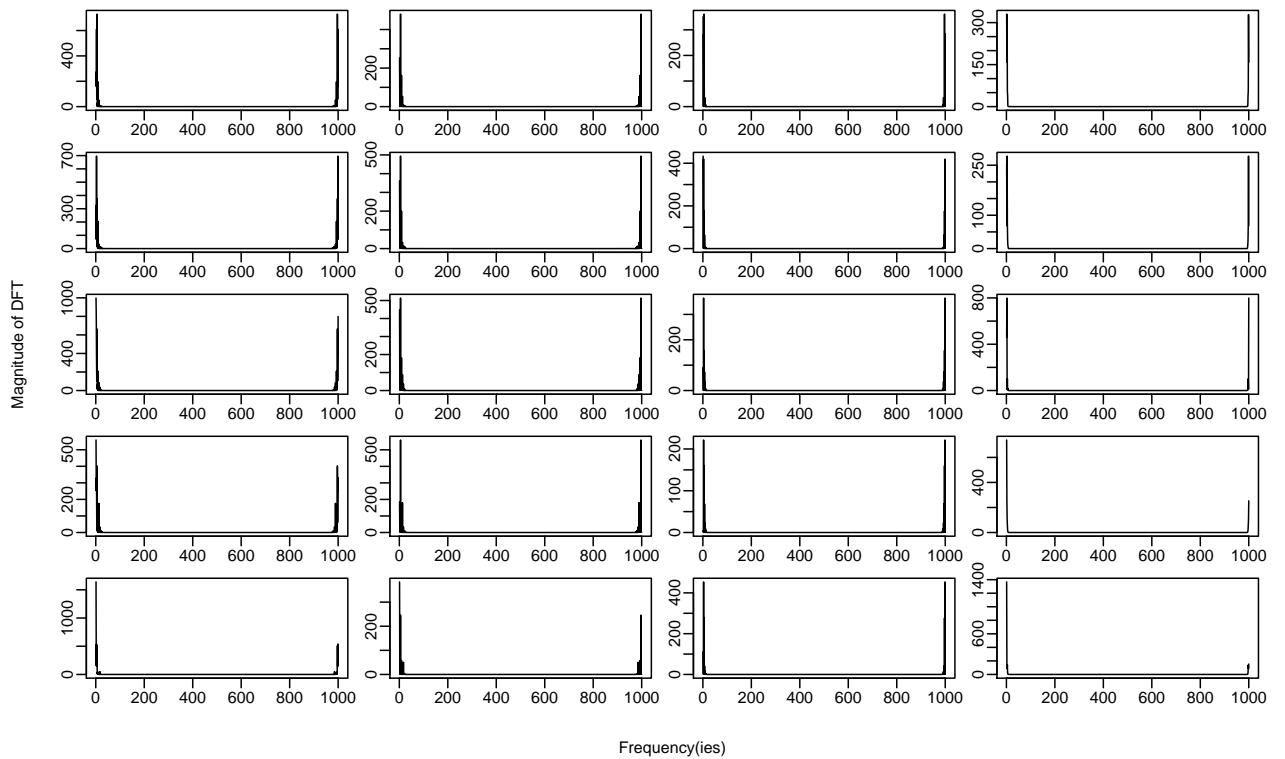


Figure 26:  $N = 1000$ ,  $l = 1$ ,  $p = 0.25, 0.5, 1.00$ . Magnitudes of the DFT,  $m = 1, 2, \dots, 10$ .

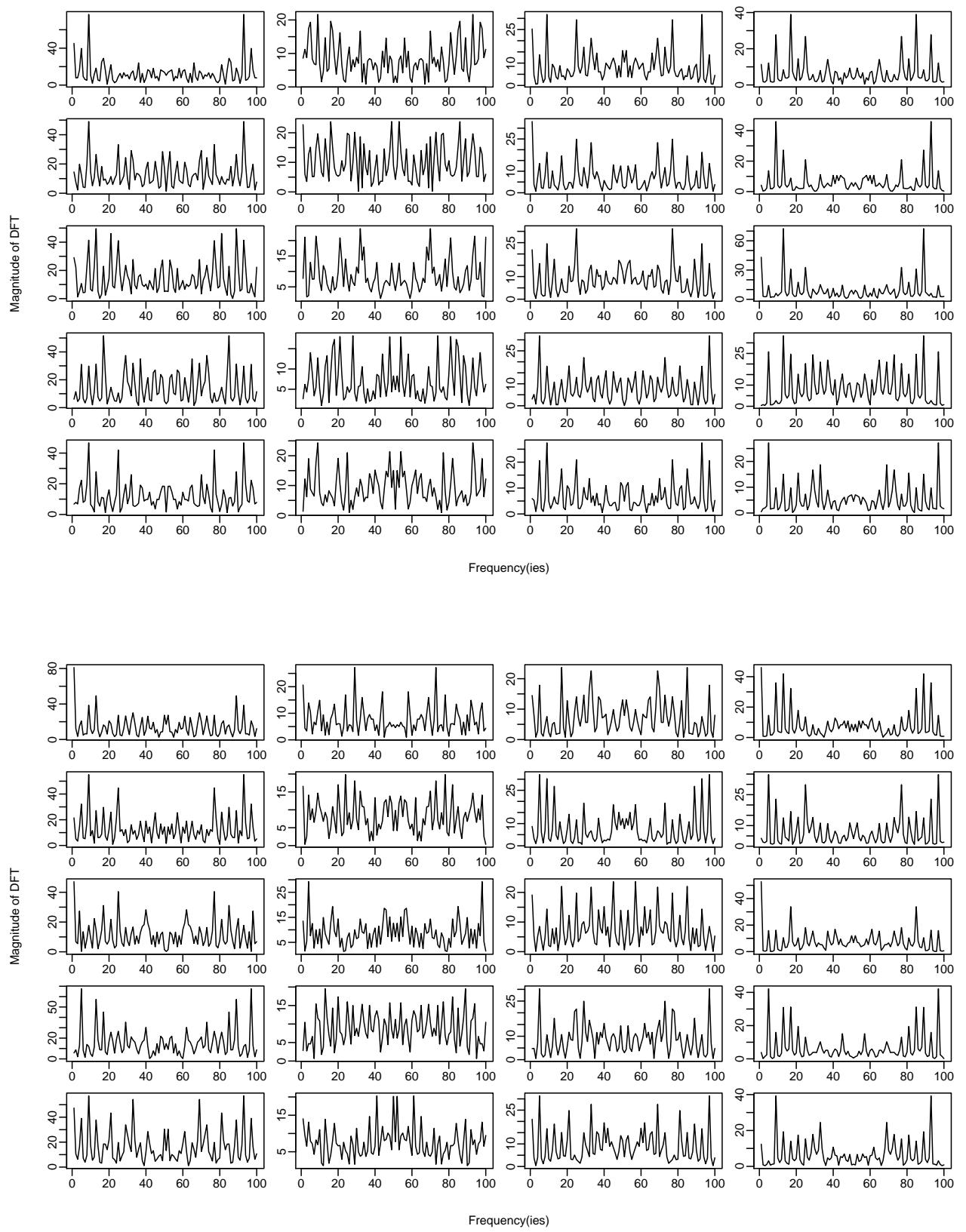


Figure 27:  $N = 100$ ,  $l = 0.05, 0.1, 0.15$ ,  $p = 0.25$ . Magnitudes of the DFT,  $m = 1, 2, \dots, 10$ .

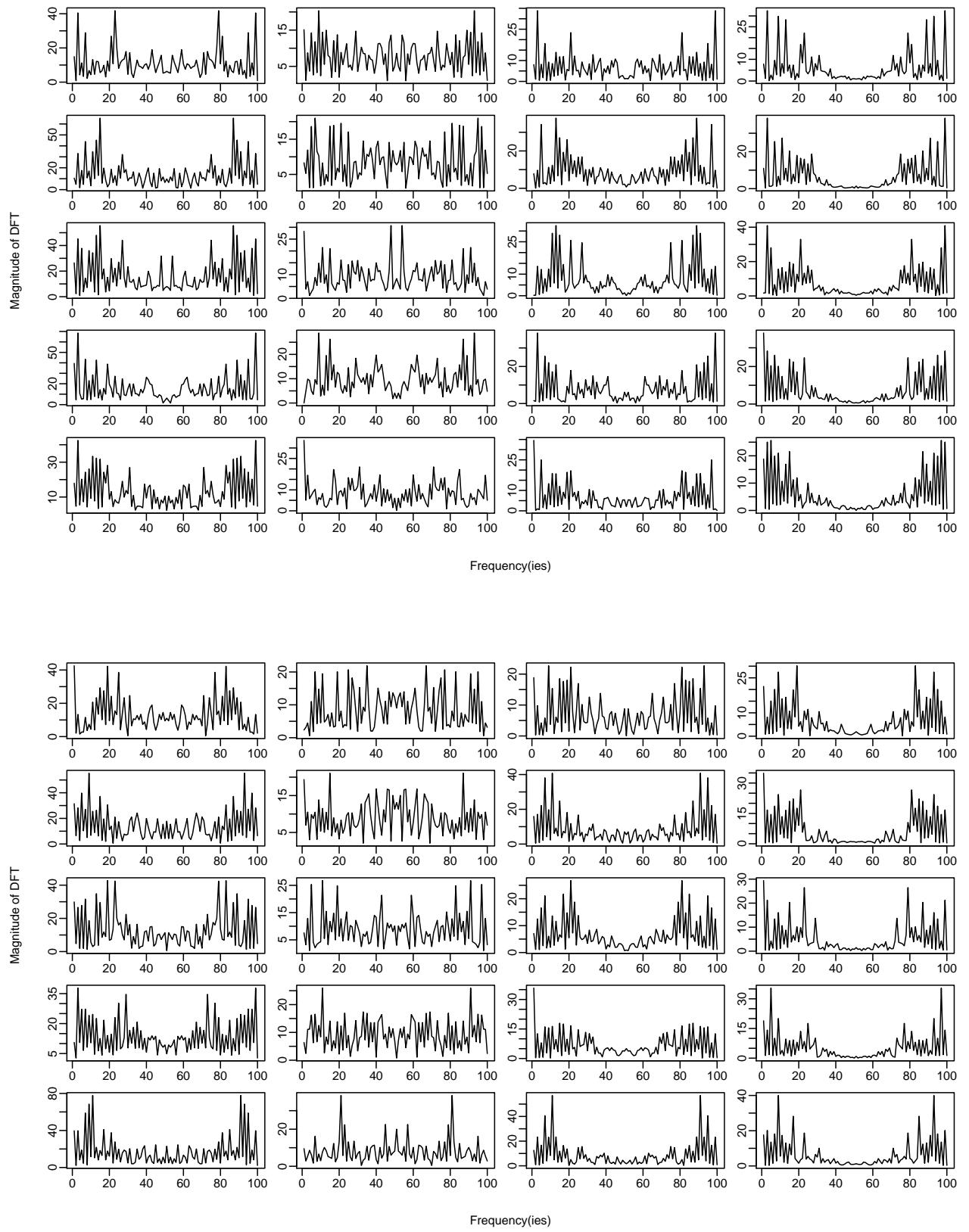


Figure 28:  $N = 100$ ,  $l = 0.05, 0.1, 0.15$ ,  $p = 0.5$ . Magnitudes of the DFT,  $m = 1, 2, \dots, 10$ .

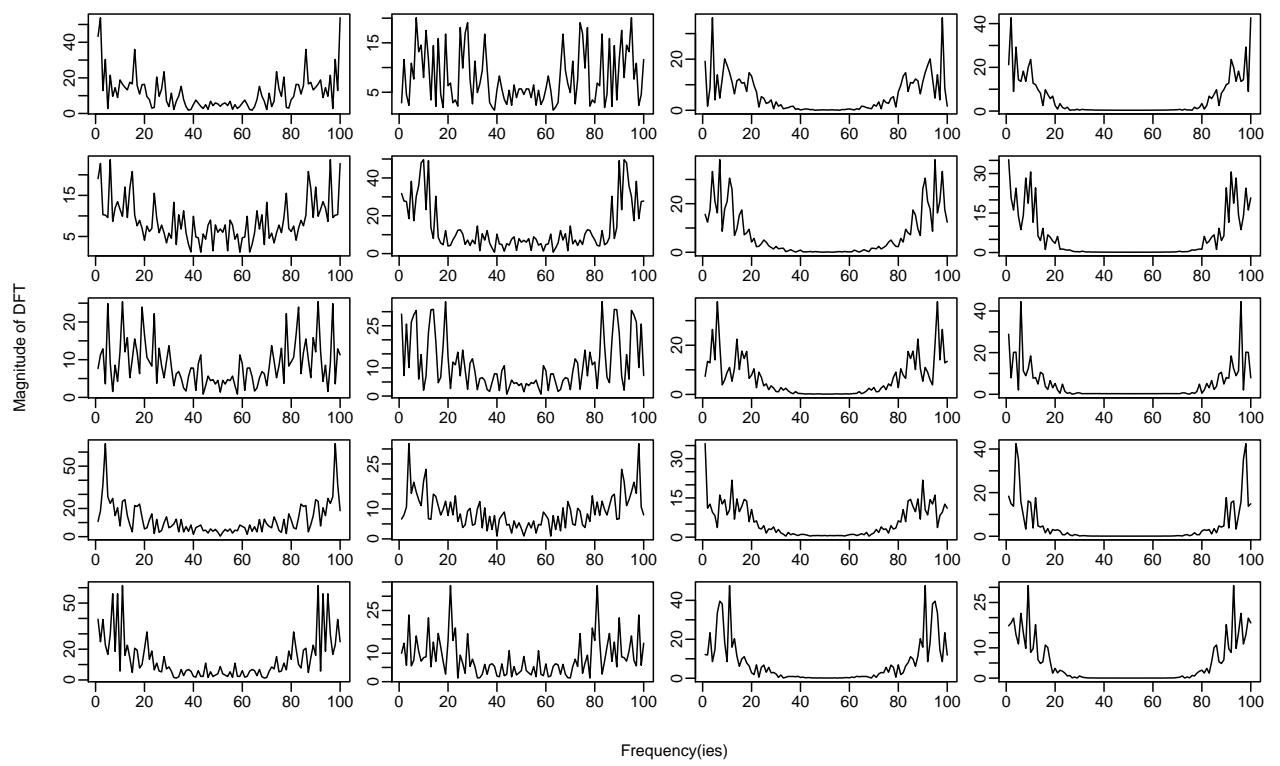
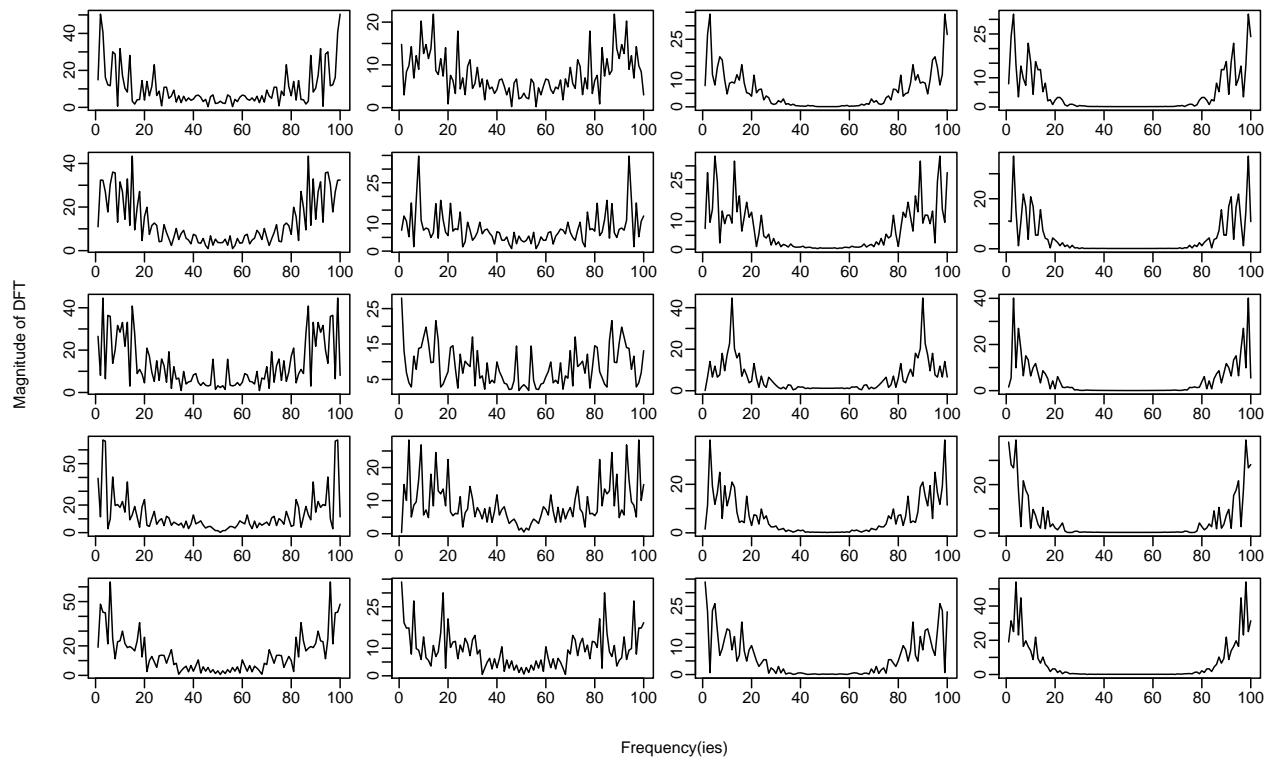


Figure 29:  $N = 100$ ,  $l = 0.05, 0.1, 0.15$ ,  $p = 1$ . Magnitudes of the DFT,  $m = 1, 2, \dots, 10$ .