

A brief description of the spectrometric data

This dataset is a part of the original one which can be found at <http://lib.stat.cmu.edu/datasets/tecolor>. For each unit i (among 215 pieces of finely chopped meat), we observe one spectrometric curve (\mathbf{x}_i) which corresponds to the absorbance measured at 100 wavelengths (i.e. $\mathbf{x}_i = (\chi_i(\lambda_1), \dots, \chi_i(\lambda_{100}))$). Moreover, for each unit i , we have at hand its fat content y_i obtained by an analytical chemical processing. The file “npfda-spectrometric.dat” contains the pairs $(\mathbf{x}_i, y_i)_{i=1, \dots, 215}$ and is organized as follows:

	Col 1	...	Col j	...	Col 100	Col 101
Row 1	$\chi_1(\lambda_1)$...	$\chi_1(\lambda_j)$...	$\chi_1(\lambda_{100})$	y_1
\vdots	\vdots	\vdots	\vdots	\vdots	\vdots	\vdots
Row i	$\chi_i(\lambda_1)$...	$\chi_i(\lambda_j)$...	$\chi_i(\lambda_{100})$	y_i
\vdots	\vdots	\vdots	\vdots	\vdots	\vdots	\vdots
Row 215	$\chi_{215}(\lambda_1)$...	$\chi_{215}(\lambda_j)$...	$\chi_{215}(\lambda_{100})$	y_{215}

The first 100 columns correspond to the 100 channel spectrum whereas the last column contains the responses.