Electronic Supplementary Information

Spectroscopic and theoretical studies on intramolecular OH---π hydrogen bonding in 4-substituted 2-allylphenols

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Supplementary Material for Organic & Biomolecular Chemistry

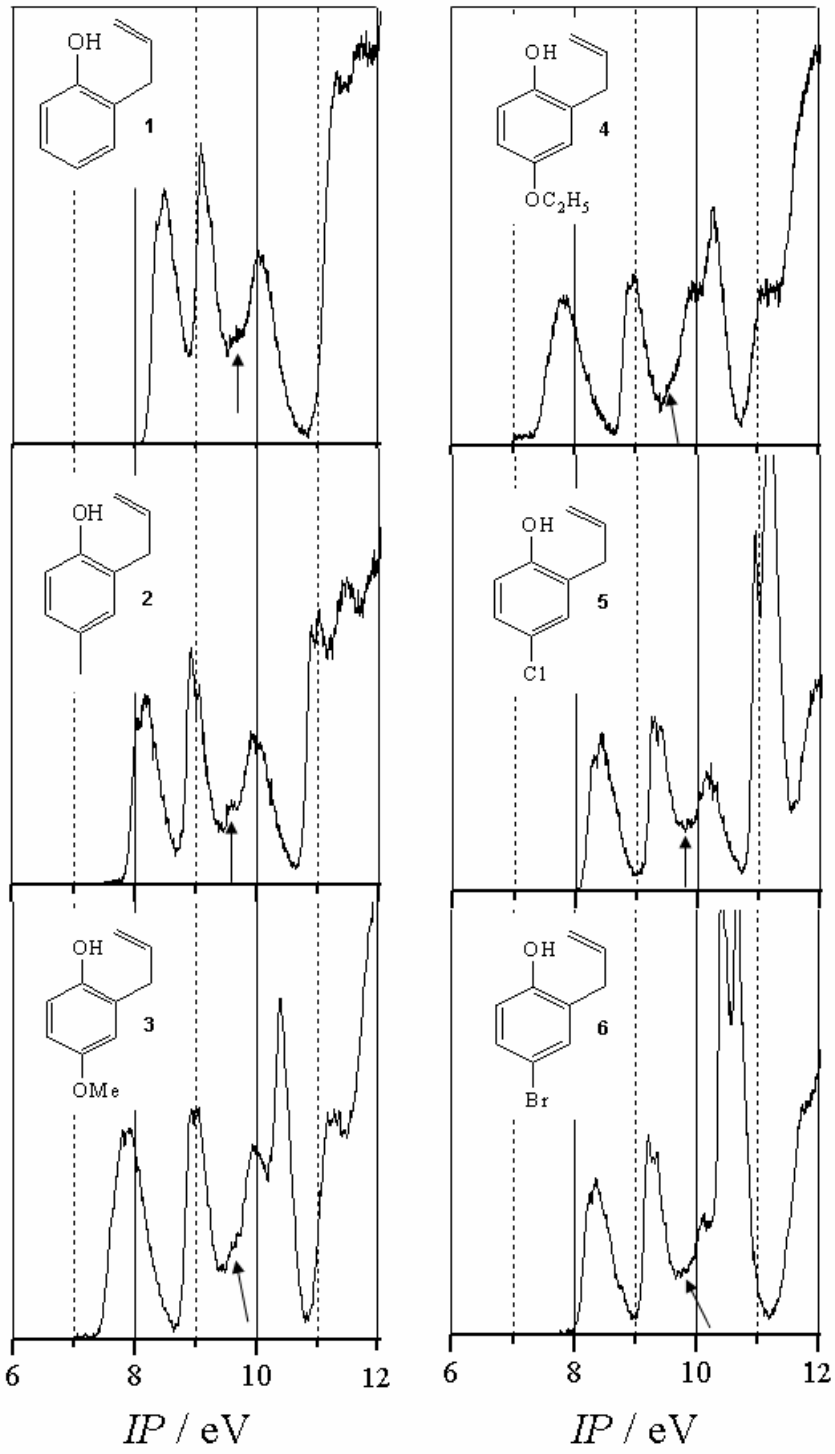

Figure ESI-1. Low-energy section of the PE spectra of 4-substituted 2-allylphenols **1**–**6**.

The approximate location of the weak ionization corresponding to conformation **b** is indicated by an arrow.

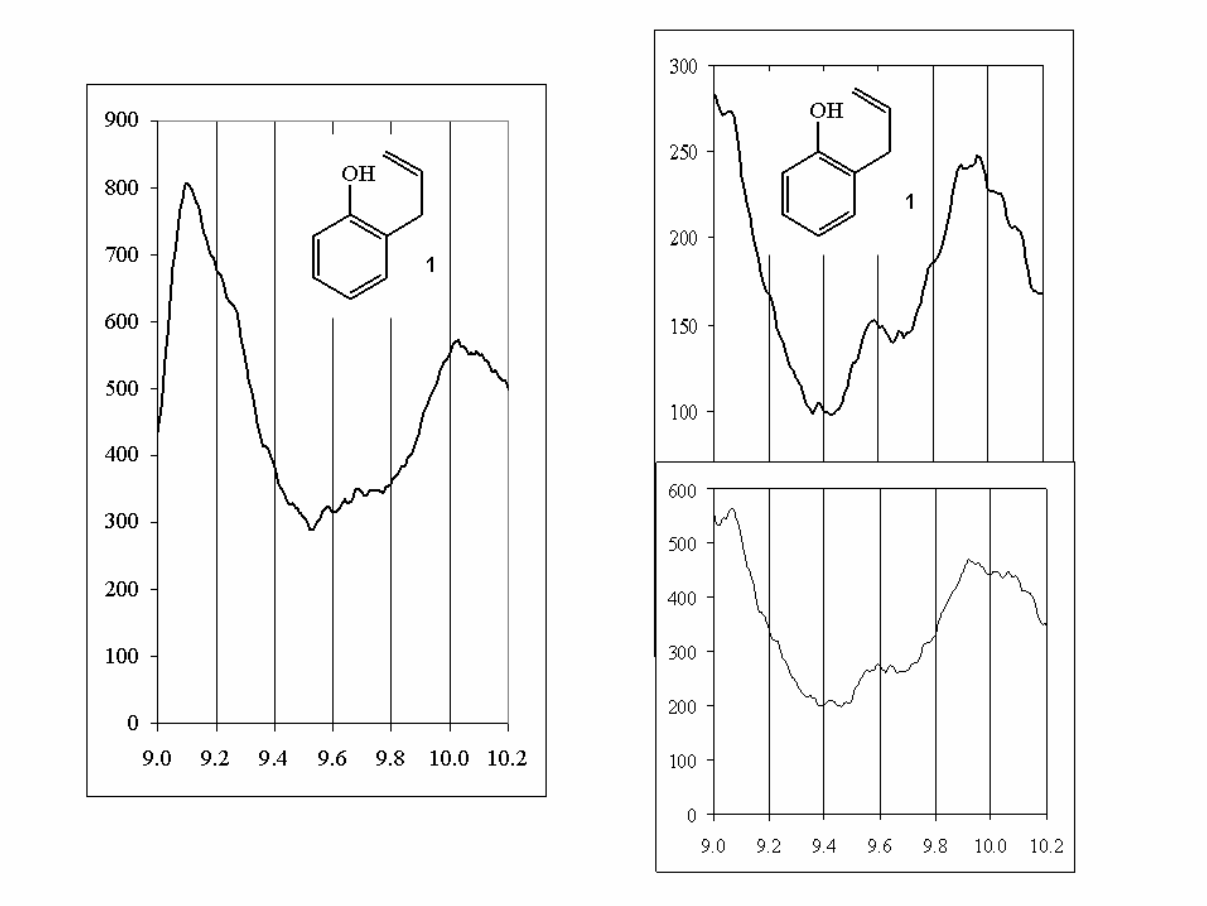

Figure ESI-3. Section of PE spectra of compound **1**.

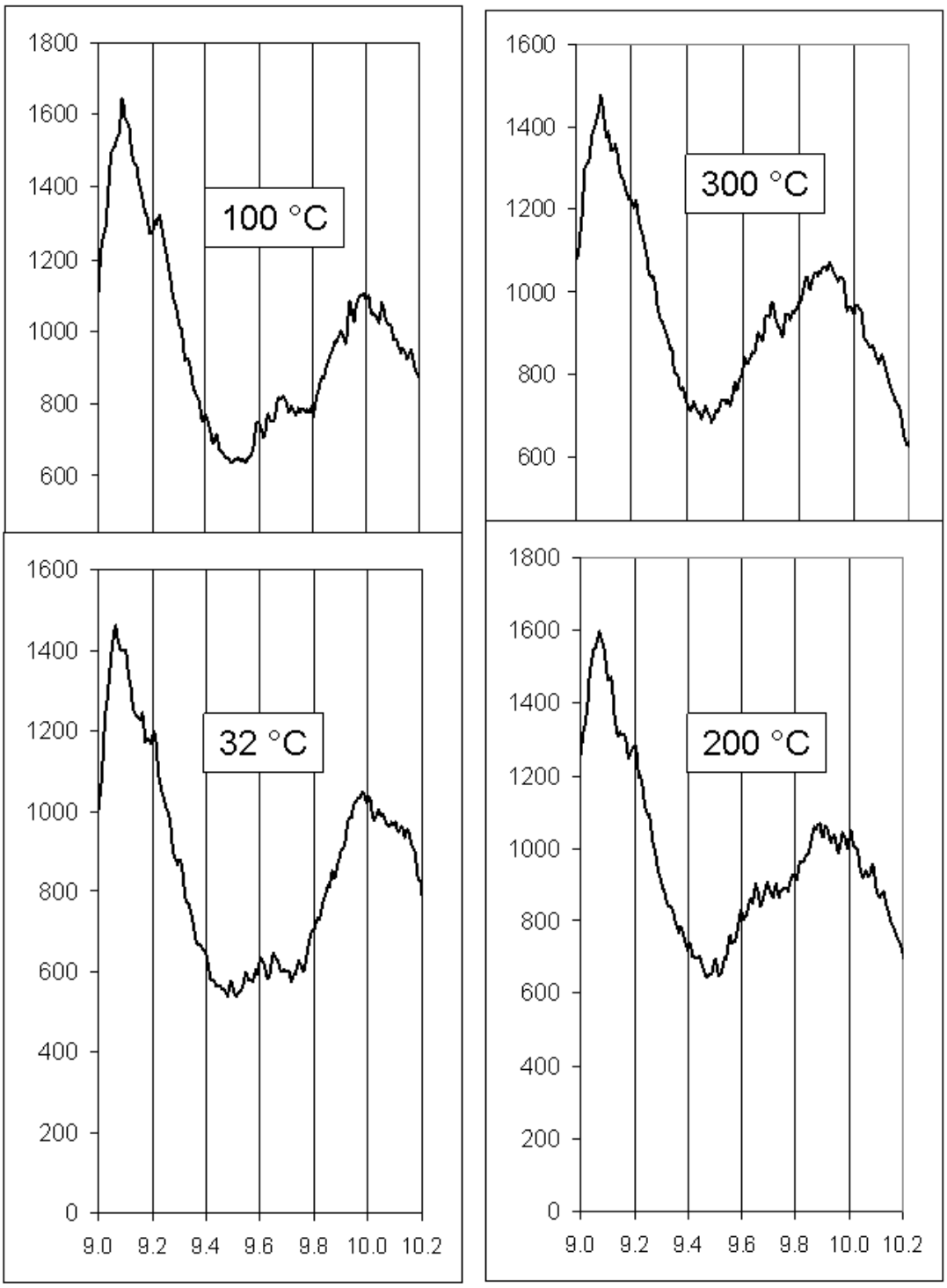

Figure ESI-3. Section of PE spectra of compound **1** recorded at different temperatures.

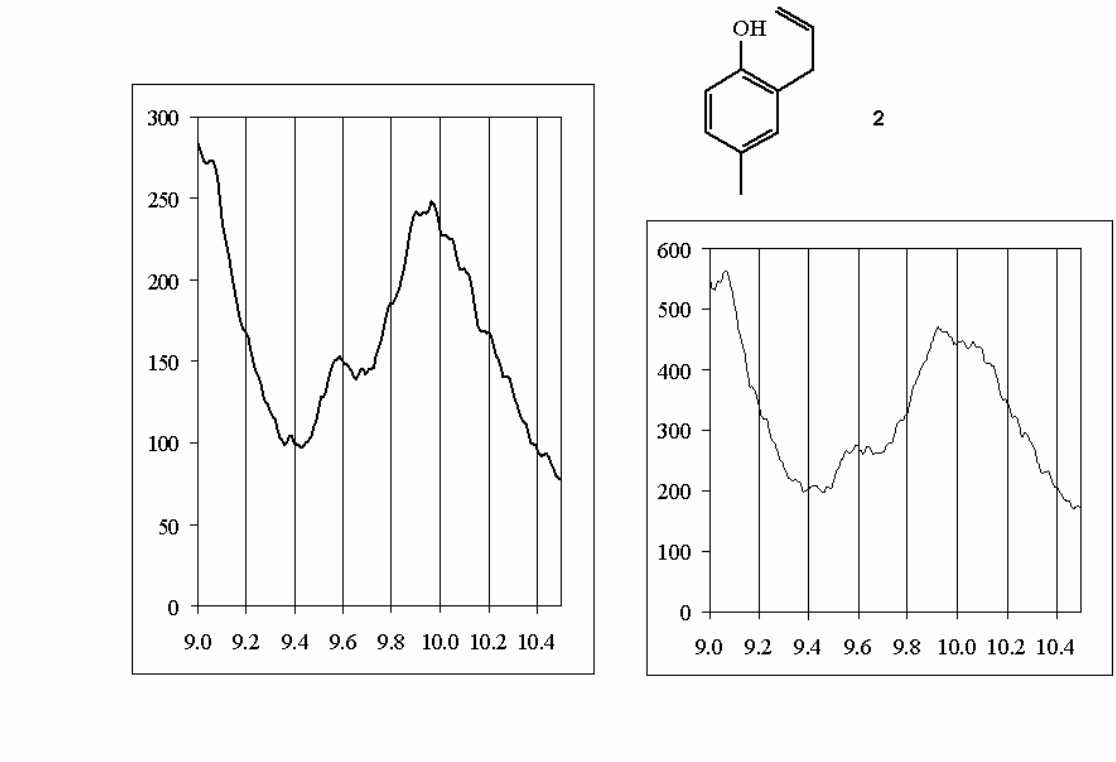

Figure ESI-4. Section of PE spectra of compound **2**.

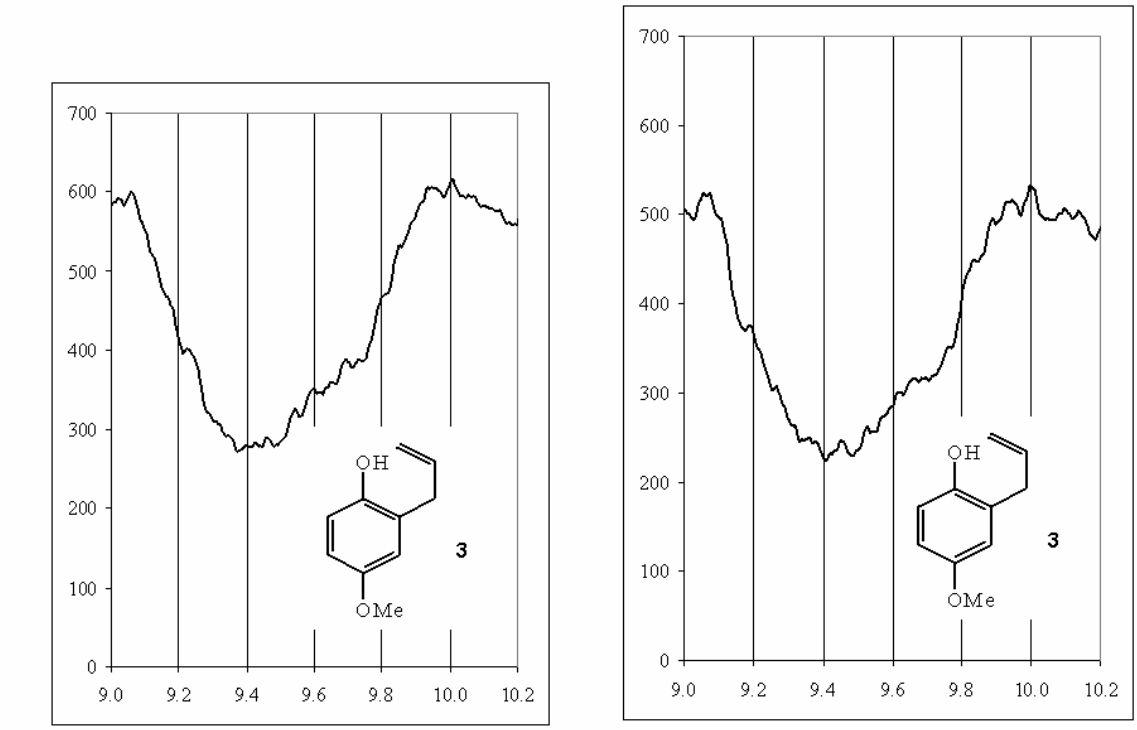

Figure ESI-5. Section of PE spectra of compound **3**.

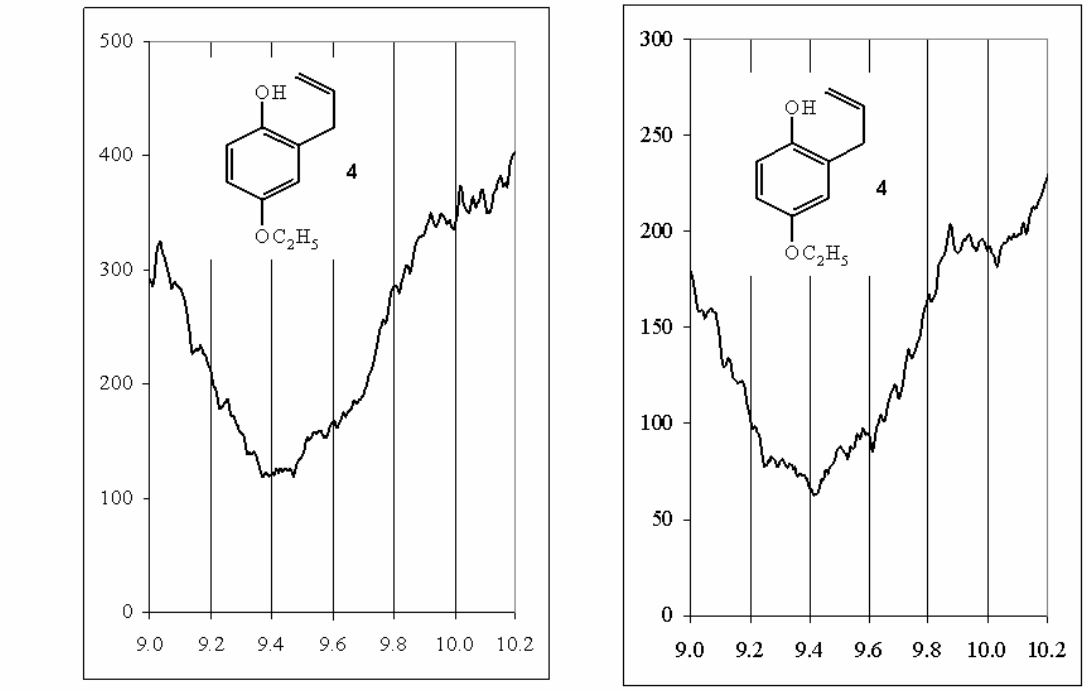

Figure ESI-6. Section of PE spectra of compound **4**.

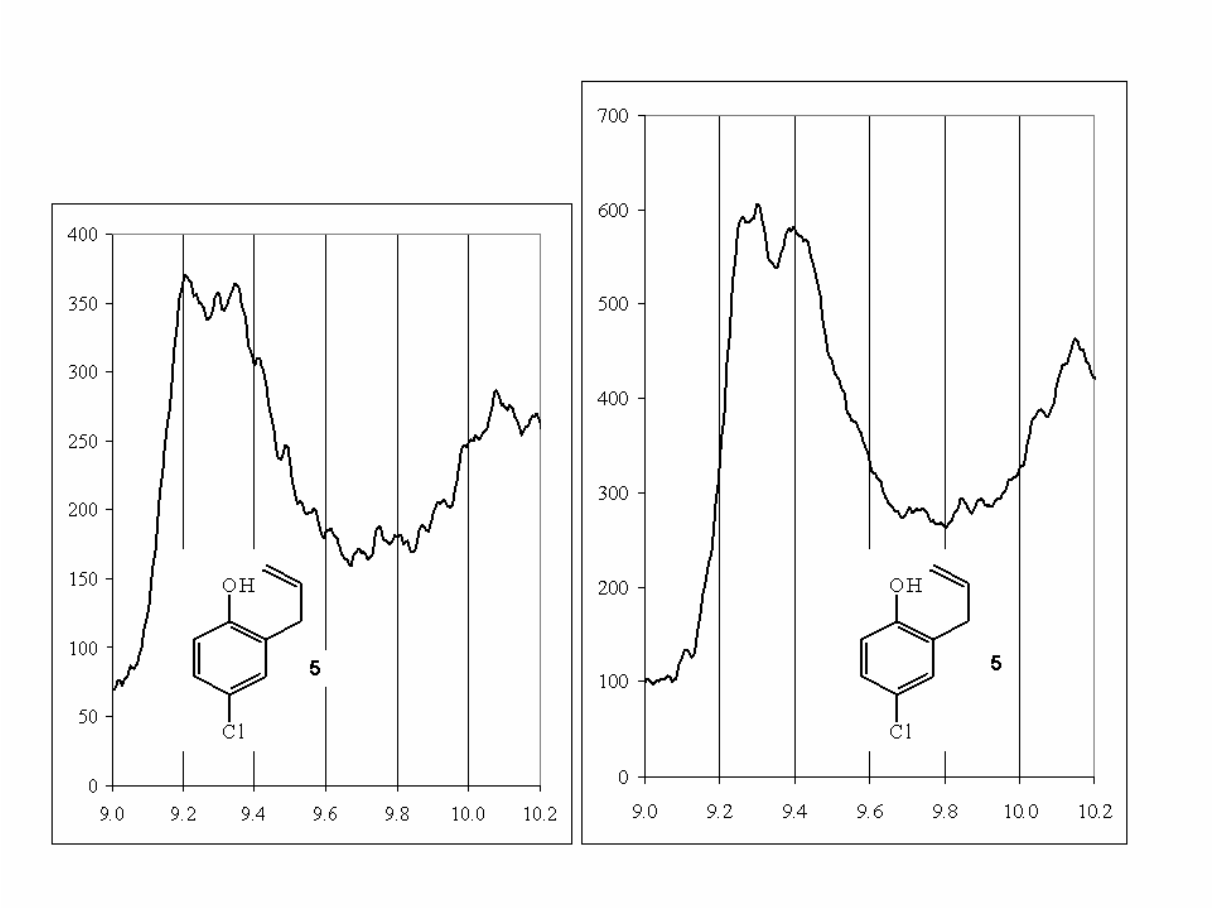

Figure ESI-7. Section of PE spectra of compound **5**.

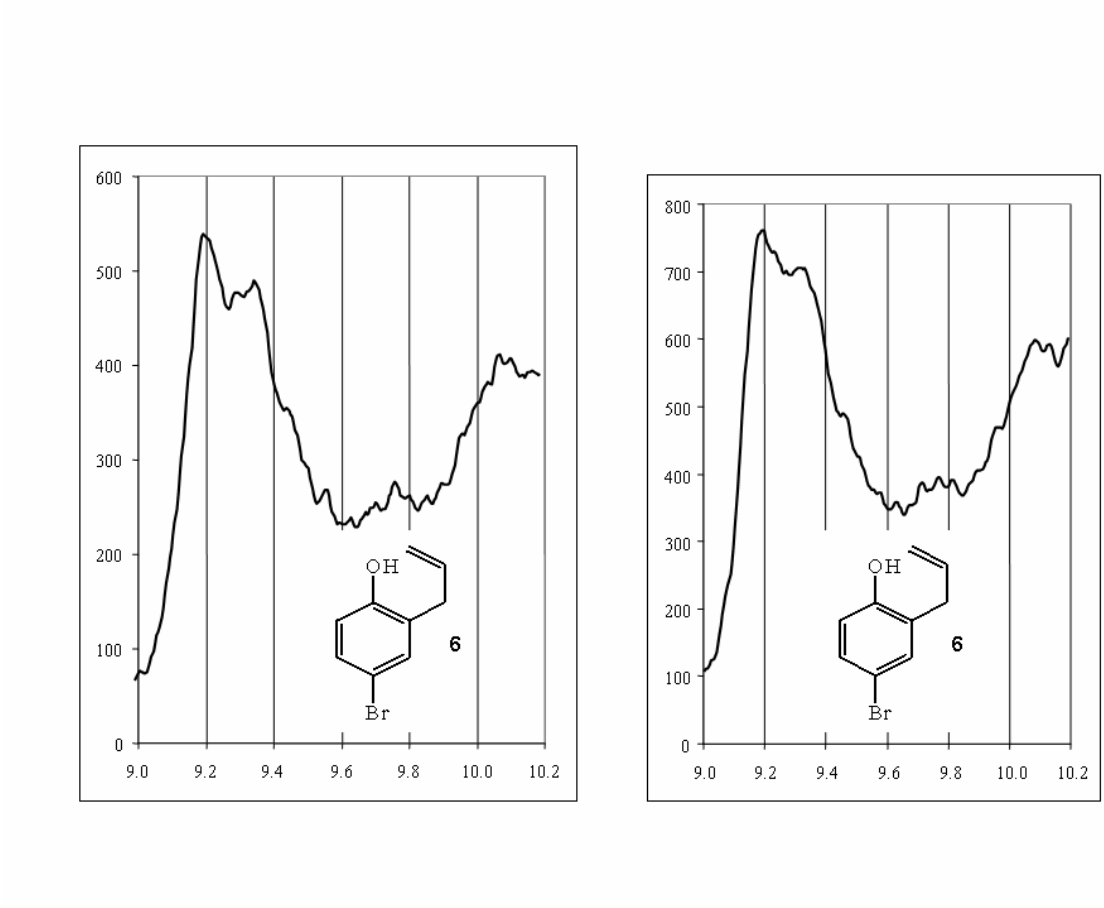

Figure ESI-8. Section of PE spectra of compound **6**.