

**Copper(II), cobalt(II), manganese(II) and nickel(II)  
bis(hexafluoroacetylacetonate) complexes with *N*-vinylimidazole**

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IR spectra were recorded on a Bruker Vertex 70 spectrometer in KBr pellets. UV–VIS spectra were recorded on a Perkin–Elmer Lambda 35 spectrophotometer. 1-Vinylimidazole, acetone, and diethyl ether were used as purchased.

*Synthesis of complexes 1–4:* 1-vinylimidazole (66 mg, 0.70 mmol) was added to a stirred solution of the corresponding hexafluoroacetylacetonate (0.30 mmol),  $[M(\text{hfac})_2(\text{H}_2\text{O})_2]$  ( $M = \text{Cu, Ni, Mn}$ ) or  $[\text{Co}(\text{hfac})_2(\text{H}_2\text{O})_2] \cdot \text{H}_2\text{O}$ , in acetone (2–3 ml). The resulting solution was filtered, and diethyl ether (5 ml) was then layered onto a filtrate. The mixture was kept overnight at room temperature. The precipitated crystals of compounds **1–4** were collected, washed with diethyl ether (1×2 ml), and dried on air.

*Complex 1* (green prisms). (72% yield). IR ( $\text{cm}^{-1}$ ): 3187, 3172, 3144, 3113, 3070 and 3020 s. ( $\nu_{\text{C-H}}$ ), 1656 s. ( $\nu_{\text{C=C}}$ , vinyl group), 1551, 1529, 1517, 1501 and 1496 s. ( $\nu_{\text{C=C}}$ ,  $\nu_{\text{C=N}}$ , imidazole) 1256, 1227, 1193, 1156 and 1130 s. ( $\nu_{\text{C-F}}$ ,  $\text{CF}_3$ ). Anal. Calcd for  $\text{C}_{20}\text{H}_{14}\text{CuF}_{12}\text{N}_4\text{O}_4$ : C, 36.07; H, 2.12; N, 8.41. Found: C, 35.92; H, 2.24; N, 8.25.

*Complex 2* (orange prisms). (80% yield). IR ( $\text{cm}^{-1}$ ): 3181, 3160, 3139, 3112, 3069 and 3021 s. ( $\nu_{\text{C-H}}$ ), 1656 и 1643 s. ( $\nu_{\text{C=C}}$ , vinyl group), 1555, 1527, 1532, 1514 and 1501 s. ( $\nu_{\text{C=C}}$ ,  $\nu_{\text{C=N}}$ , imidazole), 1255, 1228, 1193, 1154 and 1134 s. ( $\nu_{\text{C-F}}$ ,  $\text{CF}_3$ ). Anal. Calcd for  $\text{C}_{20}\text{H}_{14}\text{CoF}_{12}\text{N}_4\text{O}_4$ : C, 36.33; H, 2.13; N, 8.47. Found: C, 36.40; H, 2.31; N, 8.29.

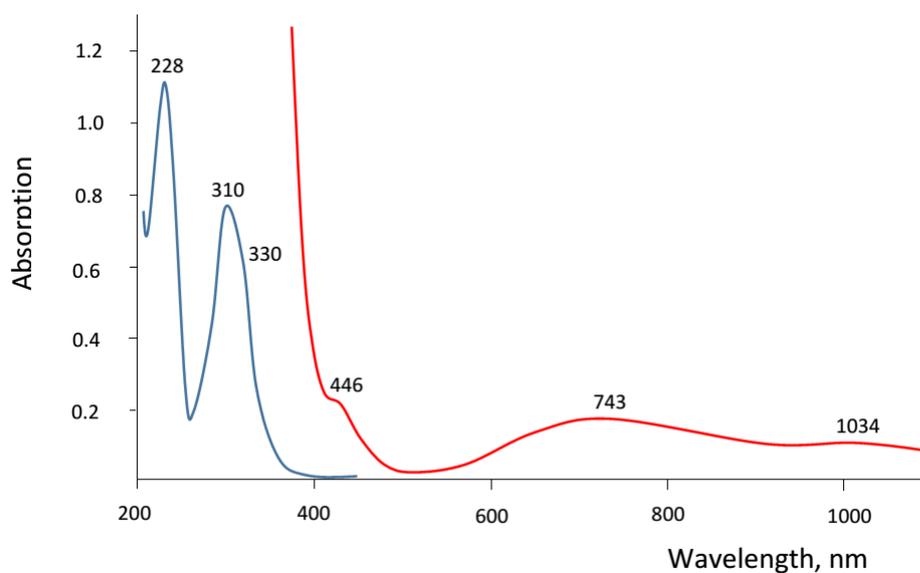
*Complex 3* (yellow prisms). (78% yield). IR ( $\text{cm}^{-1}$ ): 3180, 3153, 3133, 3110, 3066 and 3021 s. ( $\nu_{\text{C-H}}$ ), 1653 and 1649 s. ( $\nu_{\text{C=C}}$ , vinyl group), 1556, 1530, 1513 and 1500 s. ( $\nu_{\text{C=C}}$ ,  $\nu_{\text{C=N}}$ , imidazole), 1253, 1228, 1193, 1154 and 1134 s. ( $\nu_{\text{C-F}}$ ,  $\text{CF}_3$ ). Anal. Calcd for  $\text{C}_{20}\text{H}_{14}\text{F}_{12}\text{MnN}_4\text{O}_4$ : C, 36.55; H, 2.15; N, 8.52. Found: C, 36.44; H, 2.27; N, 8.64.

*Complex 4* (blue plates) (65% yield). IR ( $\text{cm}^{-1}$ ): 3182, 3162, 3142, 3113, 3071 and 3021 s. ( $\nu_{\text{C-H}}$ ), 1656 and 1647 s. ( $\nu_{\text{C=C}}$ , vinyl group), 1554, 1534, 1527, 1515 and 1502 s. ( $\nu_{\text{C=C}}$ ,  $\nu_{\text{C=N}}$ , imidazole),

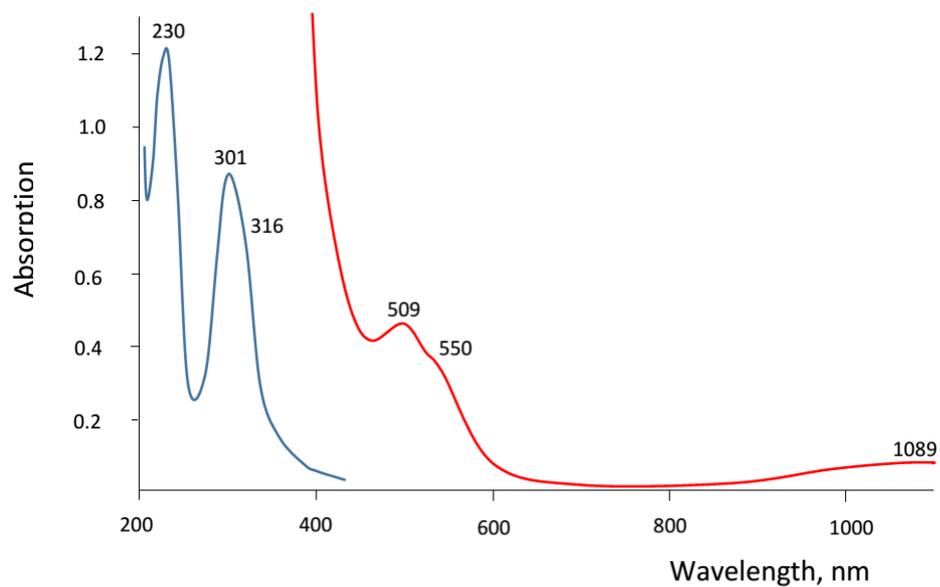
1256, 1228, 1194, 1154 and 1135 s. ( $\nu_{C-F}$ ,  $CF_3$ ). Anal. Calcd for  $C_{20}H_{14}F_{12}N_4NiO_4$ : C, 36.34; H, 2.13; N, 8.48. Found: C, 36.26; H, 2.27; N, 8.40.

**Table S1** Absorption maxima and molar extinction coefficients of complexes **1–4**.

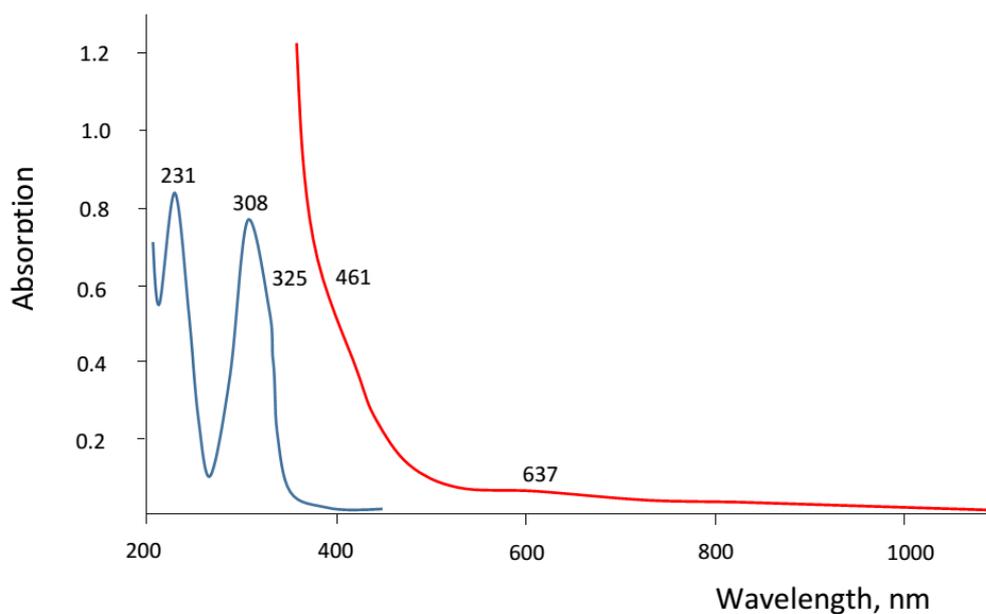
Compound	$\lambda_{max}/nm$ and $(lg\varepsilon)$						
<b>1</b>	228 (4.76)	310 (4.56)	sh 330 (4.23)	sh 446 (1.54)	743 (1.62)		sh 1034 (1.38)
<b>2</b>	230 (4.74)	301 (4.58)	sh 316 (4.58)		509 (1.79)	sh 550 (1.69)	1089 (0.84)
<b>3</b>	231 (4.77)	308 (4.77)	sh 325 (4.55)	sh 461 (1.76)	637 (1.14)		
<b>4</b>	228 (4.69)	316 (4.51)	sh 334 (4.37)	sh 438 (1.91)	609 (1.2)		1089 (0.85)



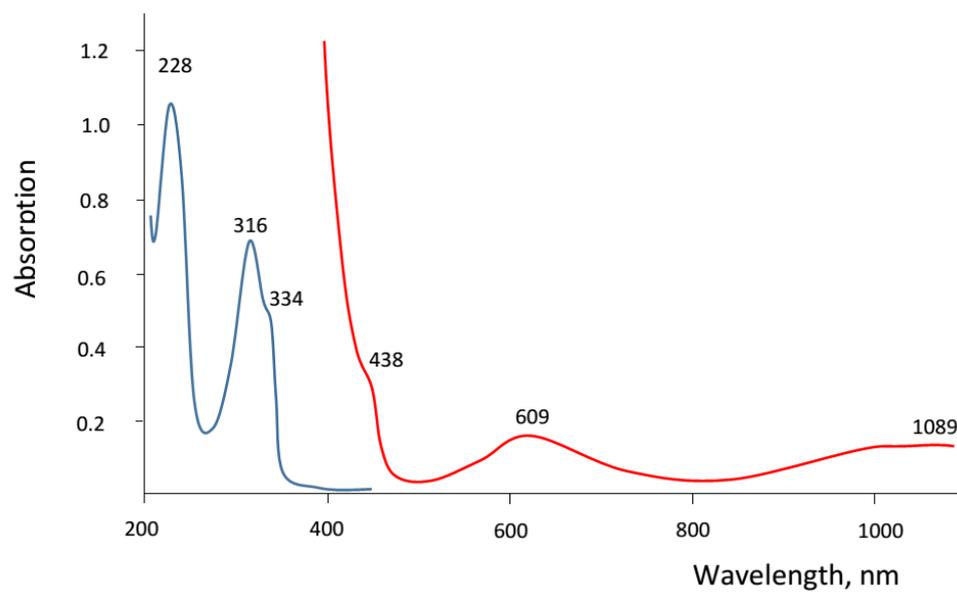
**Figure S1** Electronic absorption spectra for MeCN solutions of complex **1**:  $c = 1 \times 10^{-4}$  (blue) and  $3.9 \times 10^{-3}$  (red)  $mol\ dm^{-3}$ .



**Figure S2** Electronic absorption spectra for MeCN solutions of complex **2**:  $c = 2 \times 10^{-4}$  (blue) and  $5.4 \times 10^{-3}$  (red) mol dm<sup>-3</sup>.



**Figure S3** Electronic absorption spectra for MeCN solutions of complex **3**:  $c = 1 \times 10^{-4}$  (blue) and  $3 \times 10^{-3}$  (red) mol dm<sup>-3</sup>.



**Figure S4** Electronic absorption spectra for MeCN solutions of complex **4**:  $c = 1.5 \times 10^{-4}$  (blue)  $5.5 \times 10^{-3}$  (red) mol dm<sup>-3</sup>.