

Supplementary information

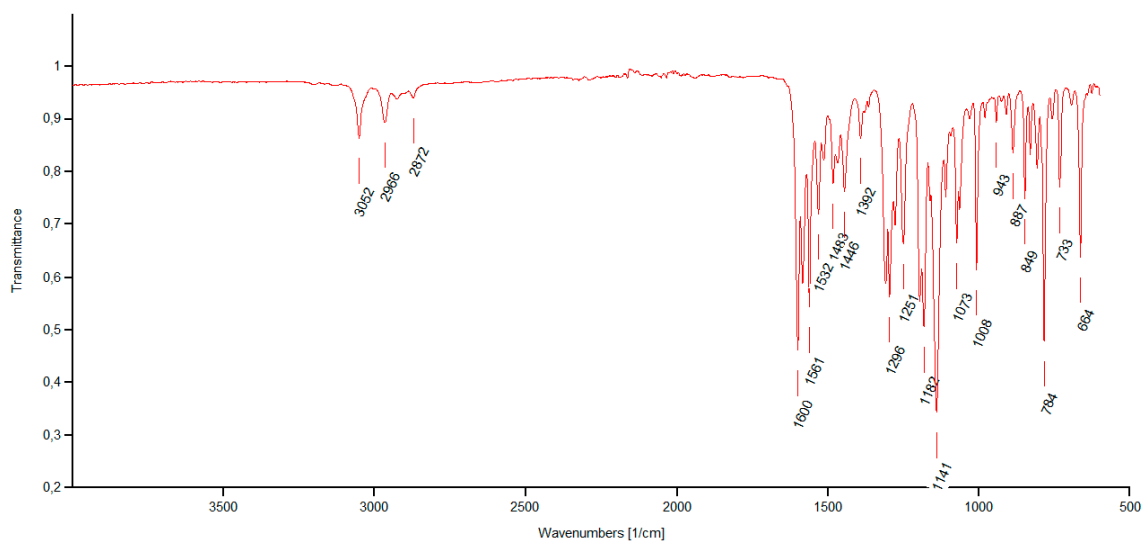


Figure 1. IR spectrum of $[(\eta^6\text{-}p\text{-cymene})\text{Ru}(4,4,4\text{-trifluoro-1-(4-bromophenyl)-1,3-butanedione})\text{Cl}]$, **1**.

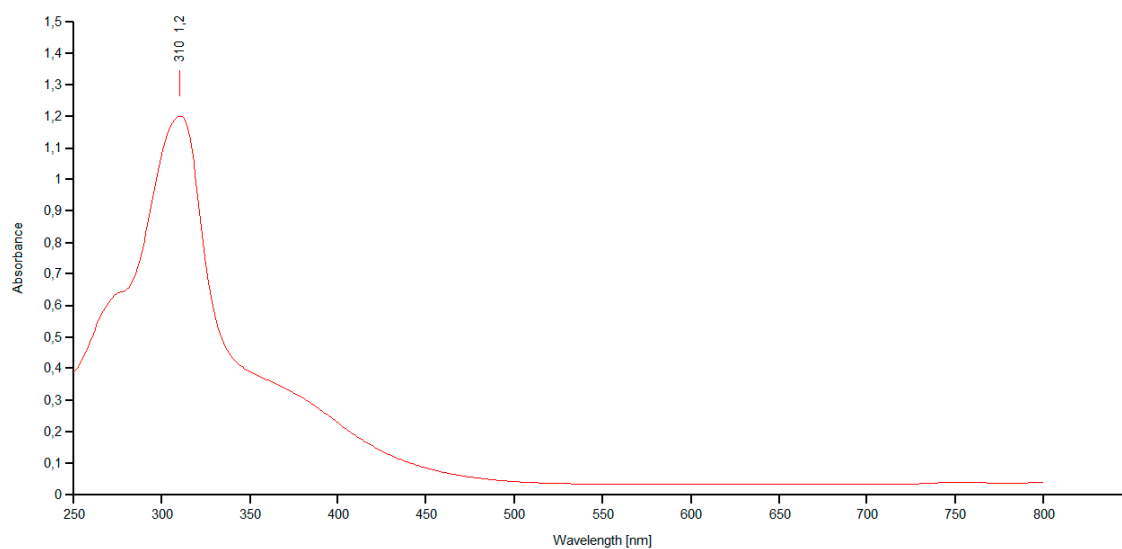


Figure 2. UV-Vis spectrum of $[(\eta^6\text{-}p\text{-cymene})\text{Ru}(4,4,4\text{-trifluoro-1-(4-bromophenyl)-1,3-butanedione})\text{Cl}]$, **1**.

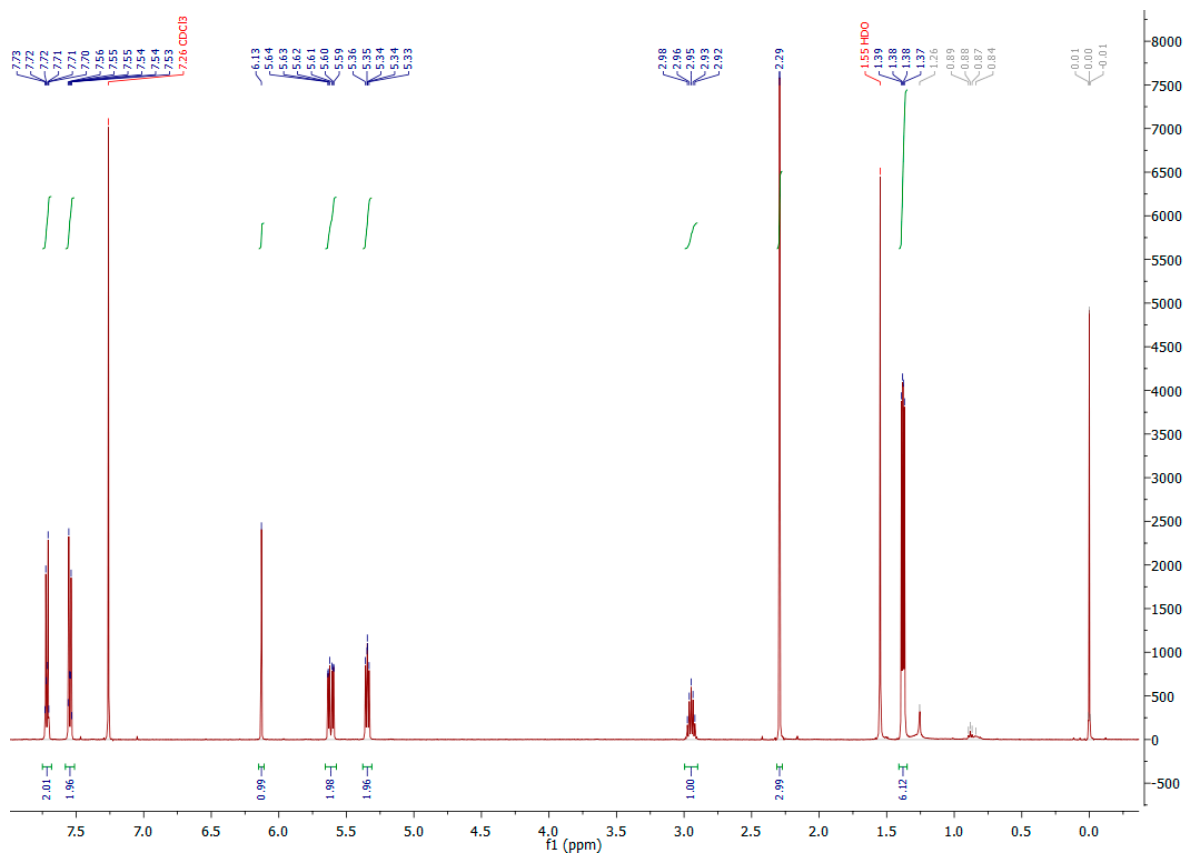


Figure 3. NMR spectrum of $[(\eta^6\text{-}p\text{-cymene})\text{Ru}(4,4,4\text{-trifluoro-1-(4-bromophenyl)-1,3-butanedione})\text{Cl}]$, **1**.

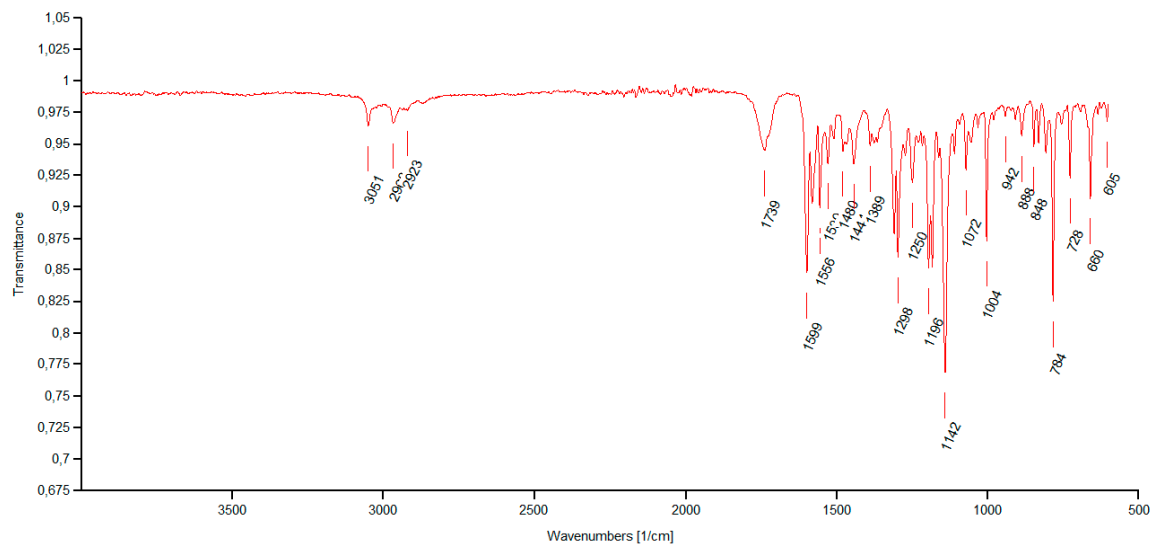


Figure 4. IR spectrum of $[(\eta^6\text{-}p\text{-cymene})\text{Ru}(4,4,4\text{-trifluoro-1-(4-iodophenyl)-1,3-butanedione})\text{Cl}]$, **3**.

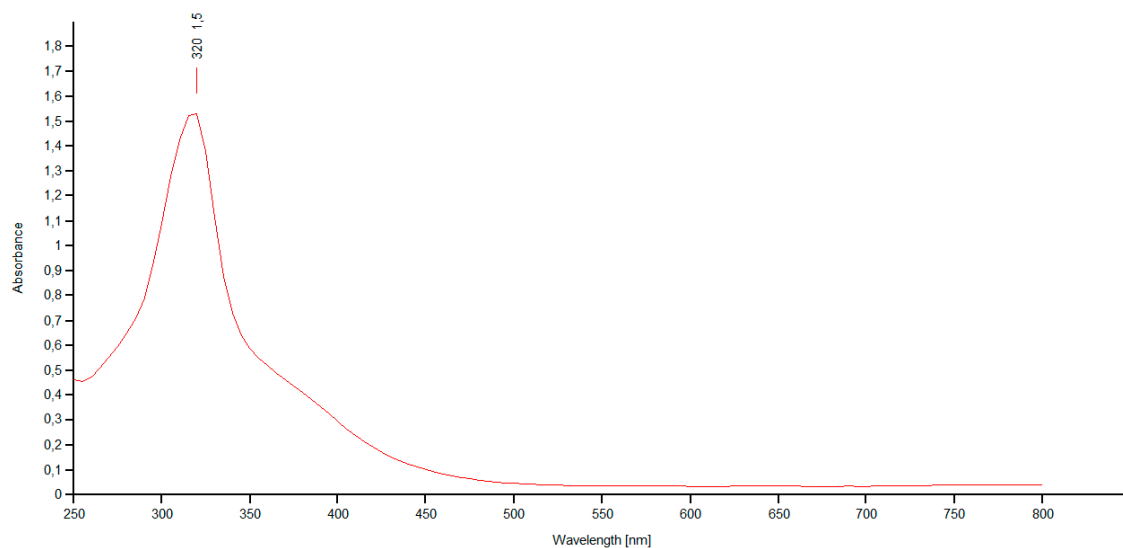


Figure 5. UV-Vis spectrum of $[(\eta^6\text{-}p\text{-cymene})\text{Ru}(4,4,4\text{-trifluoro-1-(4-iodophenyl)-1,3-butanedione})\text{Cl}]$, **3**.

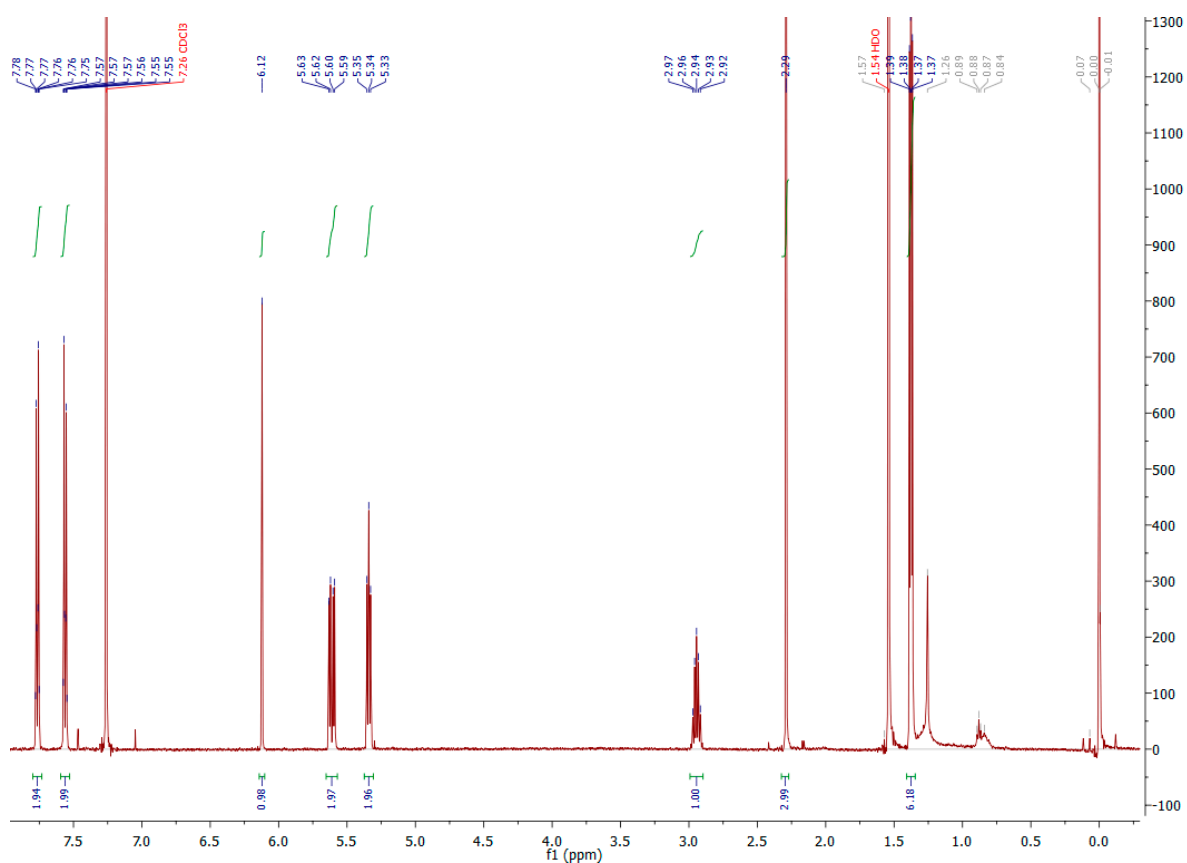


Figure 6. NMR spectrum of $[(\eta^6\text{-}p\text{-cymene})\text{Ru}(4,4,4\text{-trifluoro-1-(4-iodophenyl)-1,3-butanedione})\text{Cl}]$, **3**.

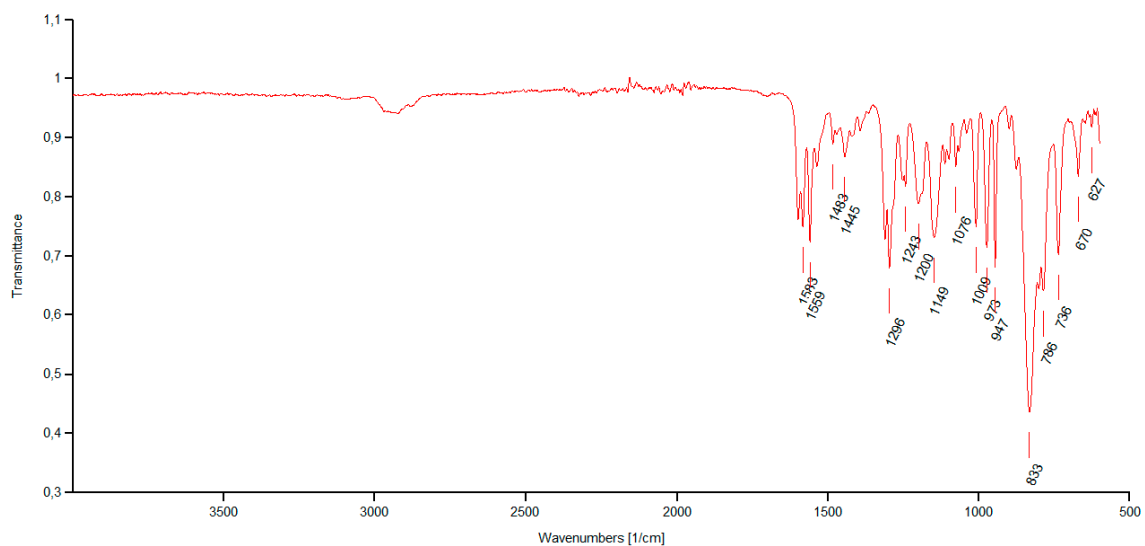


Figure 7. IR spectrum of $[(\eta^6\text{-}p\text{-cymene})\text{Ru}(4,4,4\text{-trifluoro-1-(4-bromophenyl)-1,3-butanedione})\text{pta}]\text{PF}_6$, **2**.

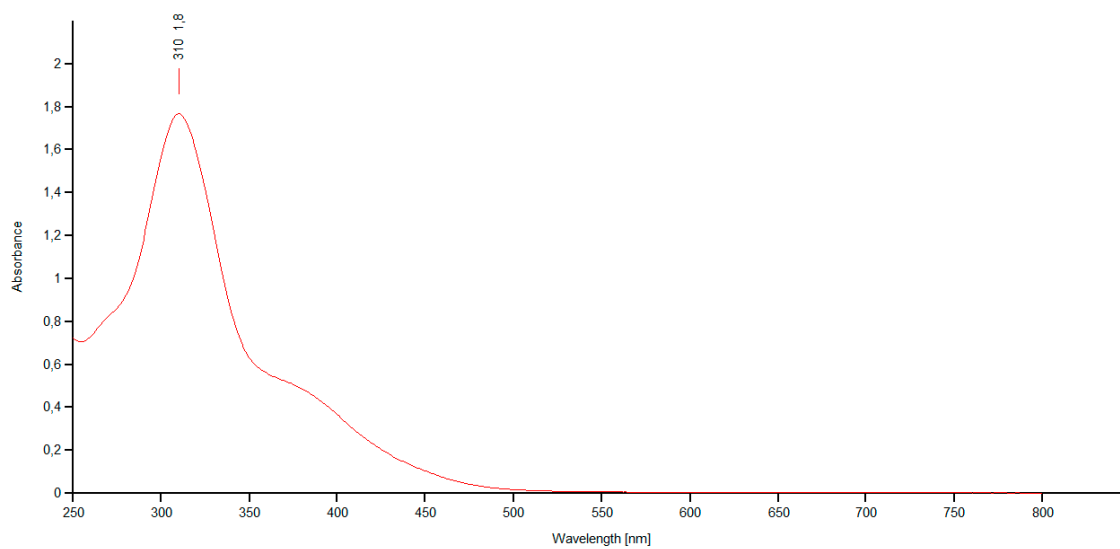


Figure 8. UV-Vis spectrum of $[(\eta^6\text{-}p\text{-cymene})\text{Ru}(4,4,4\text{-trifluoro-1-(4-bromophenyl)-1,3-butanedione})\text{pta}]\text{PF}_6$, **2**.

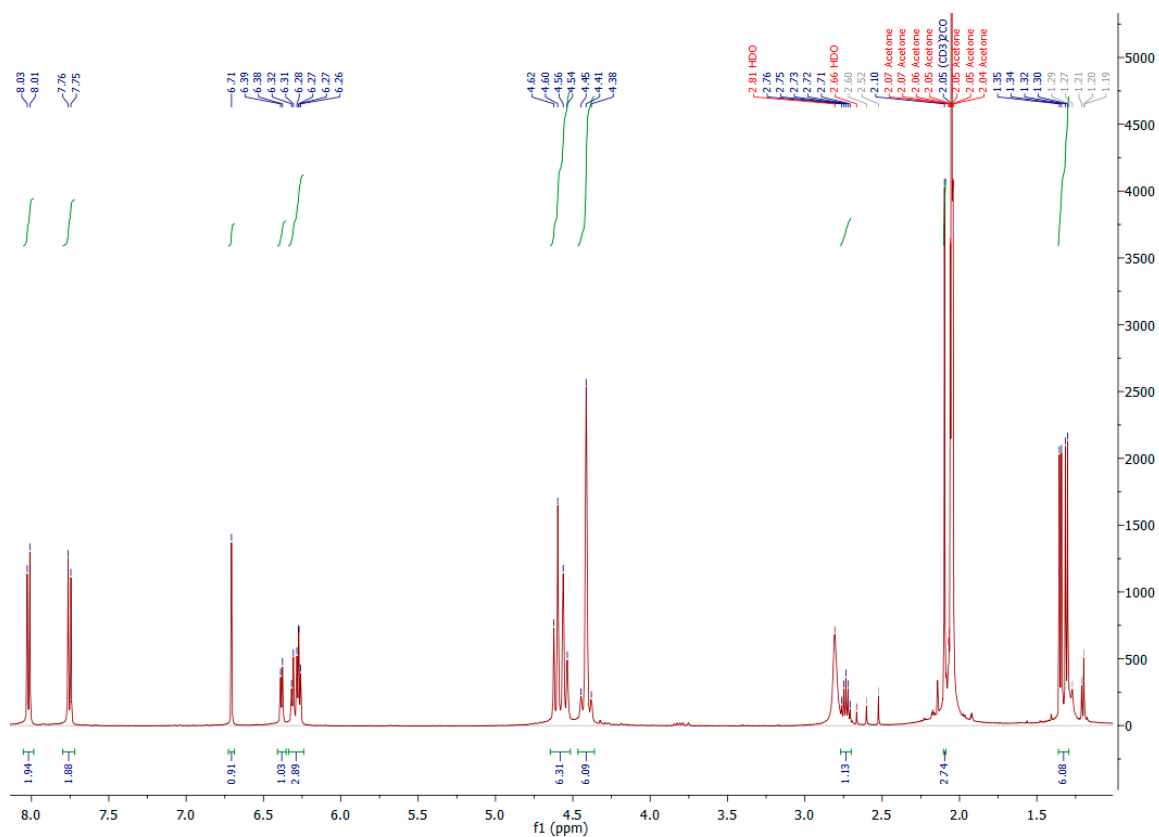


Figure 9. NMR spectrum of $[(\eta^6\text{-}p\text{-cymene})\text{Ru}(4,4,4\text{-trifluoro-1-(4-bromophenyl)-1,3-butanedione})\text{pta}]\text{PF}_6$, **2**.

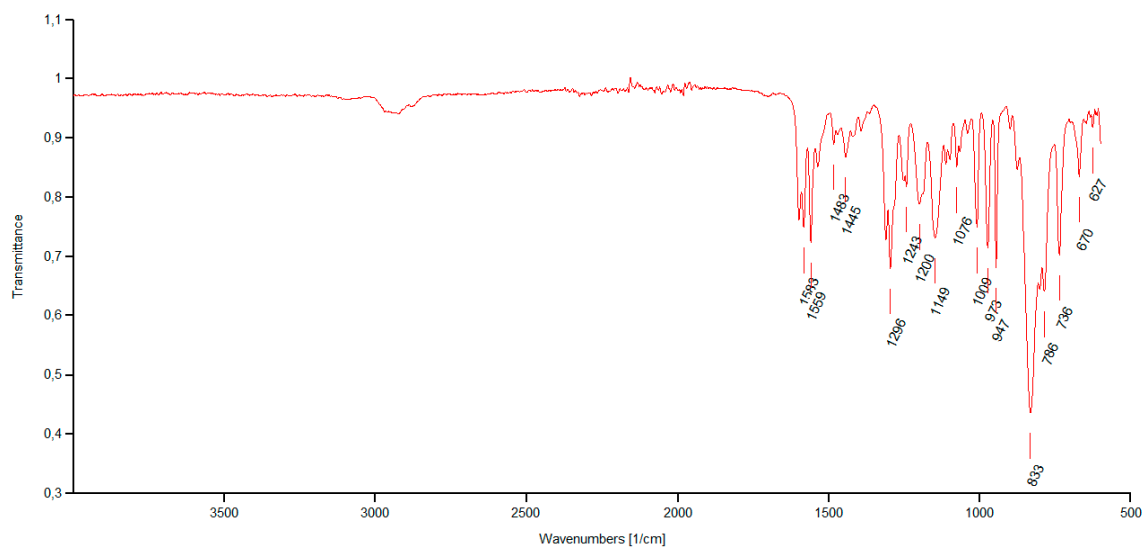


Figure 10. IR spectrum of $[(\eta^6\text{-}p\text{-cymene})\text{Ru}(4,4,4\text{-trifluoro-1-(4-iodophenyl)-1,3-butanedione})\text{pta}]\text{PF}_6$, **4**.

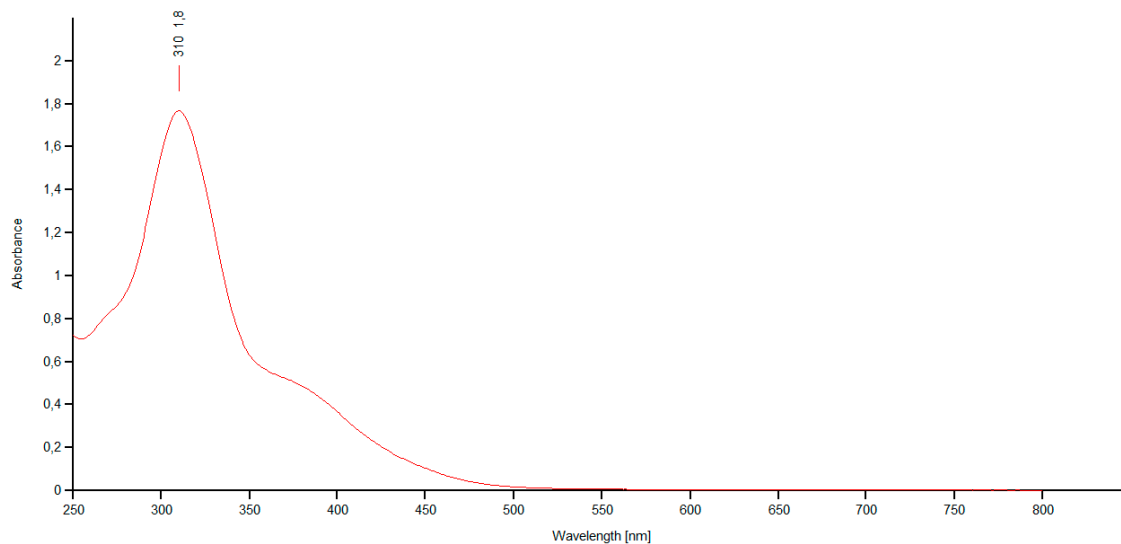


Figure 11. UV-Vis spectrum of $[(\eta^6\text{-}p\text{-cymene})\text{Ru}(4,4,4\text{-trifluoro-1-(4-iodophenyl)-1,3-butanedione})\text{pta}]\text{PF}_6$, **4**.

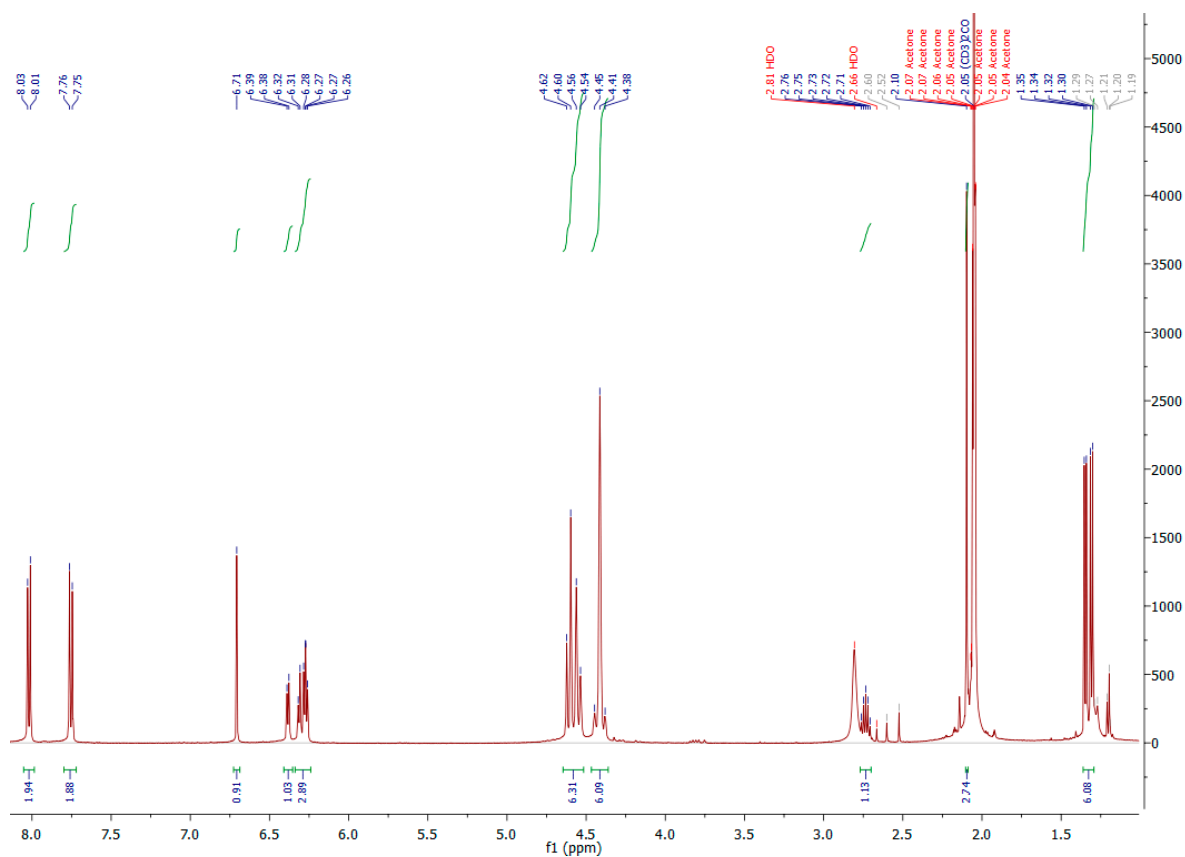


Figure 12. IR spectrum of $[(\eta^6\text{-}p\text{-cymene})\text{Ru}(4,4,4\text{-trifluoro-1-(4-iodophenyl)-1,3-butanedione})\text{pta}]\text{PF}_6$, **4**.

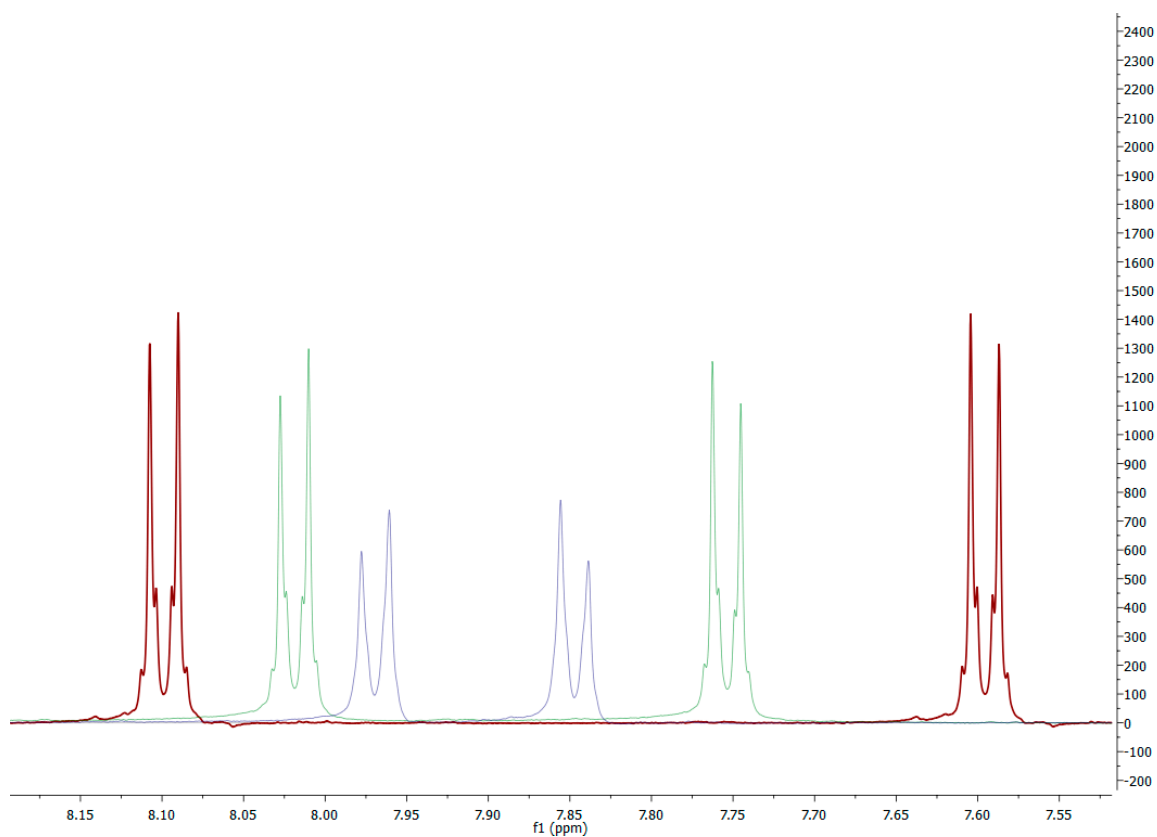


Figure 13. ^1H -NMR shifts of the hydrogen atoms on the phenyl ring of the ligand for complexes **2** (green), **4** (blue) and **6** (red).

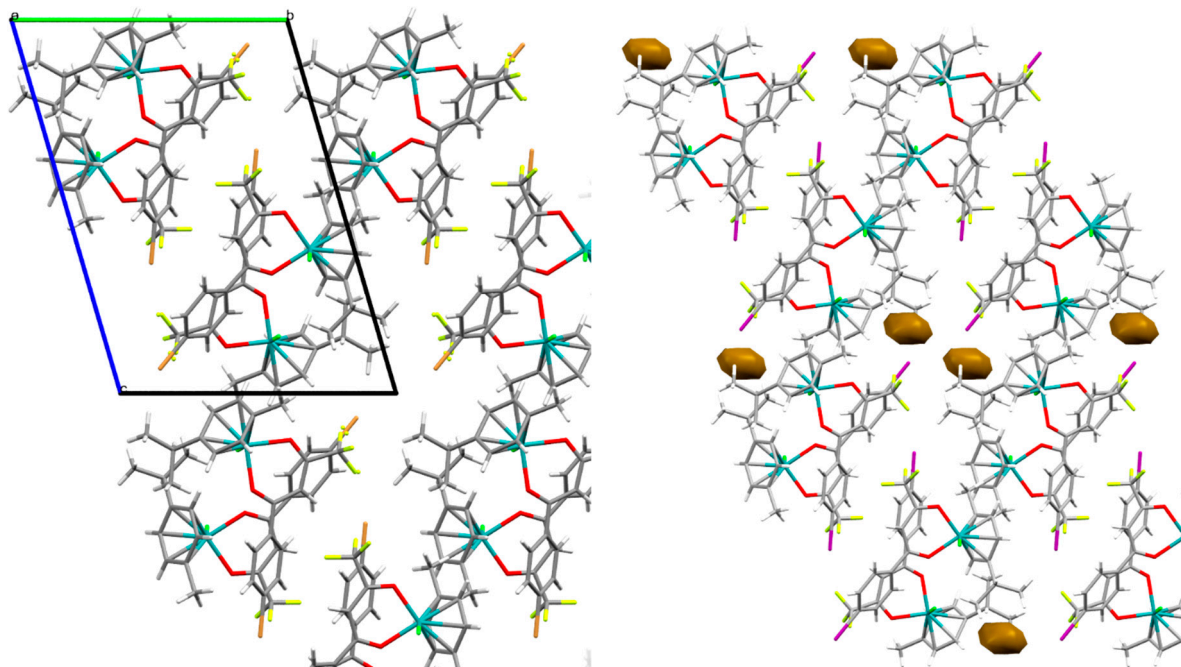


Figure 14. Packing of **1** (left) and **3** (right). Voids are indicated by ochre clouds.

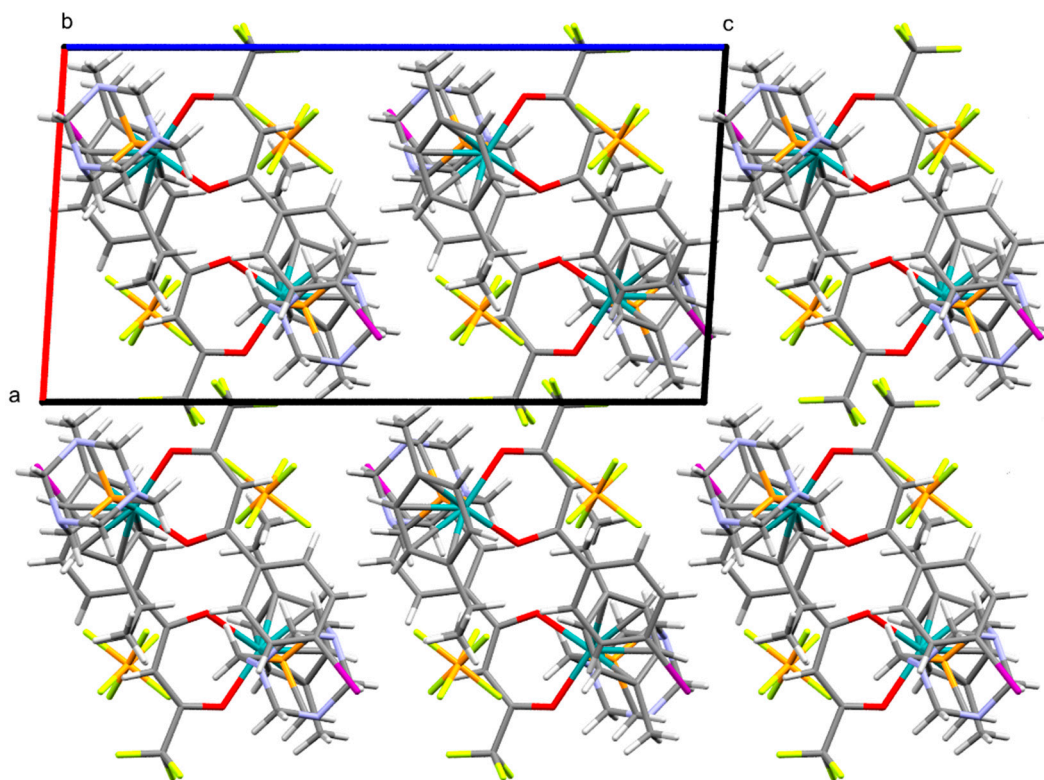


Figure 15. Packing of **4**. Note that the protruding $-\text{CF}_3$ groups are not in the same plane and do not interact with any atoms.

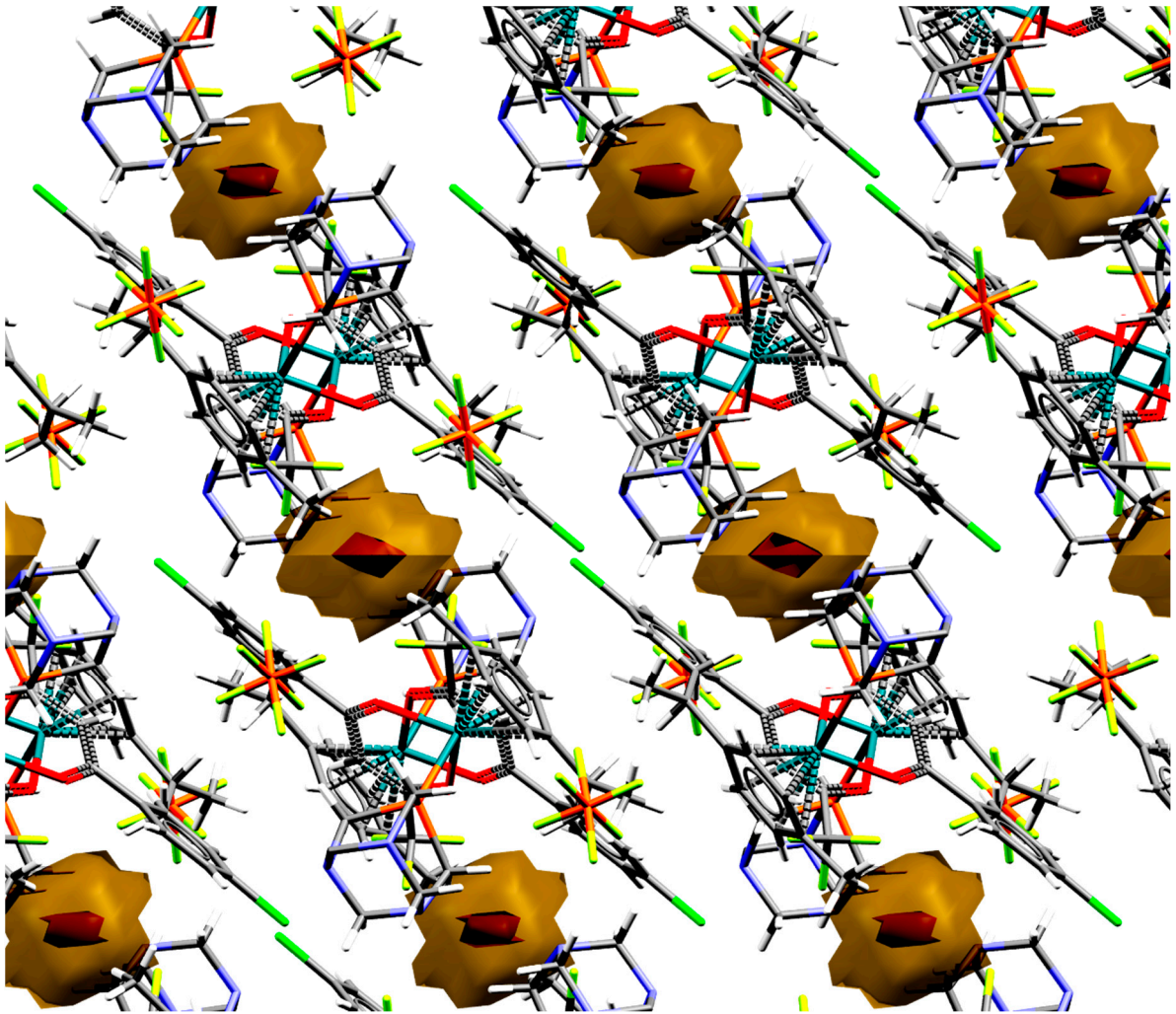


Figure 16. Packing of 6 along b axis. Solvent inaccessible voids are indicated with ochre clouds, solvent accessible voids are indicated with red clouds.