

Hydrated halogenides of Eu(II): structure, spectroscopy and charge density analysis

Autorzy

Przemysław Starynowicz

Rok wydania

2024

Czasopismo

New Journal of Chemistry

Numer woluminu

48

Strony

15921-15934

DOI

10.1039/d4nj02731k

Kolekcja

Naukowa

Język

Angielski

Typ publikacji

Artykuł

Streszczenie

Five hydrated halogenides of divalent europium, $\text{EuCl}_2 \cdot 2\text{H}_2\text{O}$ (tagged as **Cl2**), $\text{EuCl}_2 \cdot 6\text{H}_2\text{O}$ (**Cl6**), $\text{EuBr}_2 \cdot \text{H}_2\text{O}$ (**Br1**), $\text{EuBr}_2 \cdot 6\text{H}_2\text{O}$ (**Br6**) and $\text{EuI}_2 \cdot 6\text{H}_2\text{O}$ (**I6**), were studied herein. The crystals are all isomorphous with analogous compounds of Sr. In **Cl2**, the metal cation is surrounded by 4 Cl^- anions and 4 water molecules; in **Br1**, it is surrounded by 7 Br^- anions and 2 water molecules; and in **Cl6**, **Br6** and **I6**, it is surrounded by 9 water molecules, with infinite $[\text{Eu}(\text{H}_2\text{O})_6]^{2\oplus\oplus}_{\infty}$ chains being formed. Spectroscopic properties of **Cl2**, **Br1**, **Br6** and **I6** were studied. **Cl2** and **Br1** are bright luminophores, whereas **Br6** and **I6** are weak emitters at room temperature. In the latter two compounds, long-wave emissions at about 550 nm were observed and theoretical calculations indicated that this was due to $6,7s \rightarrow 4f$ transitions. Additionally, $4f \rightarrow 4f$ transitions were observed in the excitation spectra of **Br6** and **I6** recorded at 77 K. Experimental charge density analysis was performed for **Cl6**, and the results show that Eu–O bonds in the $[\text{Eu}(\text{H}_2\text{O})_6]^{2\oplus\oplus}_{\infty}$ polymeric aqua cation are more ionic than those in $[\text{Gd}(\text{H}_2\text{O})_9]^{3+}$.

Adres publiczny

<http://dx.doi.org/10.1039/d4nj02731k>

Strona internetowa wydawcy

<https://www.rsc.org/>