## Supporting Information

Effect of Substituent Groups on Crystal Structures and Luminescence Properties of 2D/3D Cd ${ }^{\text {II }}$ Complexes with Mixed Rigid and Flexible Carboxylate Ligands<br>Guo-Yun Wu, ${ }^{\text {a }}$ Yi-Xia Ren, ${ }^{\text {ab }}$ Zheng Yin, ${ }^{\text {a }}$ Feng Sun, ${ }^{\text {a Ming-Hua Zeng*a }}$ and Mohamedally Kurmoo ${ }^{c}$<br>${ }^{a}$ Key Laboratory for the Chemistry and Molecular Engineering of Medicinal Resources (Ministry of Education of China), School of Chemistry and Pharmaceutical Sciences, Guangxi Normal University, Guilin 541004, P. R. China, E-mail:zmh@mailbox.gxnu.edu.cn<br>${ }^{b}$ College of Chemistry and Chemical Engineering, Shaanxi Key Laboratory of Chemical Reaction Engineering, Yan'an University, Yan'an 716000, P. R. China<br>${ }^{c}$ Institut de Chimie de Strasbourg, Université de Strasbourg, CNRS-UMR 7177, 4 rue Blaise Pascal, CS 90032,<br>67081 Strasbourg Cedex, France, E-mail: kurmoo@inistra.fr


$l\left(\kappa^{1}\right)-\left(\kappa^{1}\right)-\mu_{2}-a$
 IV cis- $\left(\kappa^{1}-K^{1}\right)-\left(\kappa^{1}\right)-\mu_{3} \quad \vee$ trans $-\left(K^{1}-K^{1}\right)-\left(\kappa^{1}\right)-\mu_{3}$




VIII $\left(\kappa^{2}\right)-\left(\kappa^{1}\right)-\mu_{2}-b$


III $\left(\kappa^{1}\right)-\left(\kappa^{1}\right)-\mu_{2}-c$

$\mathrm{VI}\left(\kappa^{2}\right) 1$


Chart S1. The coordination modes of the Hbic ligand.


Chart S2. Reported coordination modes of adipate.


Figure S1. The 2D structure based on bic ligand in complex 1.


Figure S2. The 1D/2D structures based on adp ${ }^{2-}$ ion of complexes $\mathbf{1}(\mathrm{a}), \mathbf{2}(\mathrm{b})$ and $\mathbf{3}(\mathrm{c})$.


Figure S3. (a) The TG curves of complex 1 in flowing $\mathrm{N}_{2}$ at a heating rate of $10{ }^{\circ} \mathrm{C} / \mathrm{min}$ in the temperature range $30-900^{\circ} \mathrm{C}$.


Figure S3. (b) The TG curves of complex 2 in flowing $\mathrm{N}_{2}$ at a heating rate of $10{ }^{\circ} \mathrm{C} / \mathrm{min}$ in the temperature range $30-900^{\circ} \mathrm{C}$.


Figure S3. (c) The TG curves of complex 3 in flowing $\mathrm{N}_{2}$ at a heating rate of $10{ }^{\circ} \mathrm{C} / \mathrm{min}$ in the temperature range $30-900^{\circ} \mathrm{C}$.


Figure S4. (a) XRPD pattern of $\mathbf{1}$ compared to that simulated from single crystal data of $\mathbf{1}$.


Figure S4. (b) XRPD pattern of $\mathbf{2}$ compared to that simulated from single crystal data of $\mathbf{2}$.


Figure S4. (c) XRPD pattern of $\mathbf{3}$ compared to that simulated from single crystal data of $\mathbf{3}$.

