

Electronic supporting information

One-dimensional chains based on linear tetranuclear copper(I) units: reversible structural transformation and luminescence change

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Fig. S1. The solid-state UV-visible spectra for **1** and **2** at room temperature.

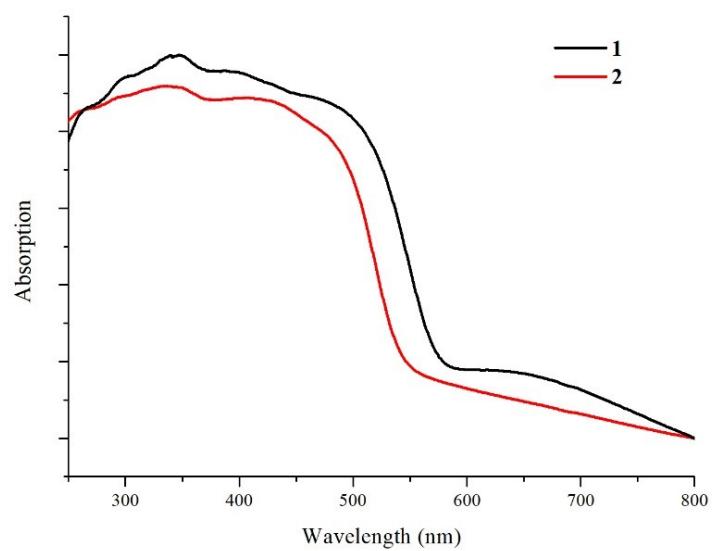


Fig. S2. The PXRD patterns of **1**. (a) Simulation and (b) synthesis.

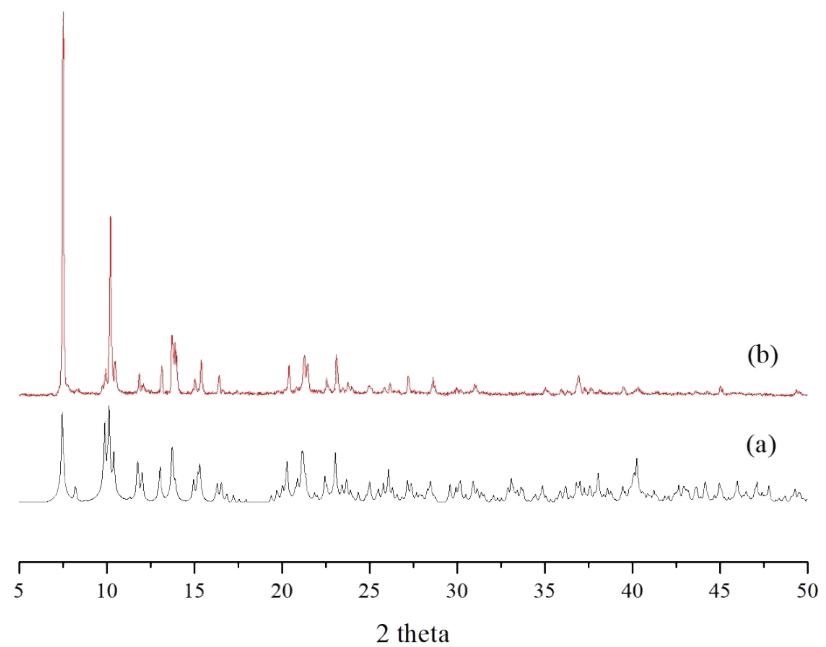


Fig. S3. The PXRD patterns of **2**. (a) Simulation and (b) synthesis.

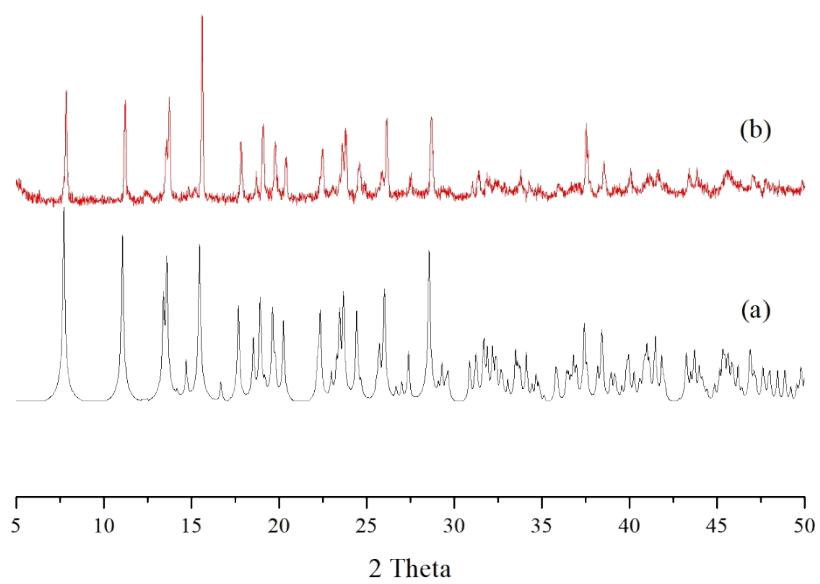


Fig. S4. PXRD patterns for complex **1** immersed in various solvents.

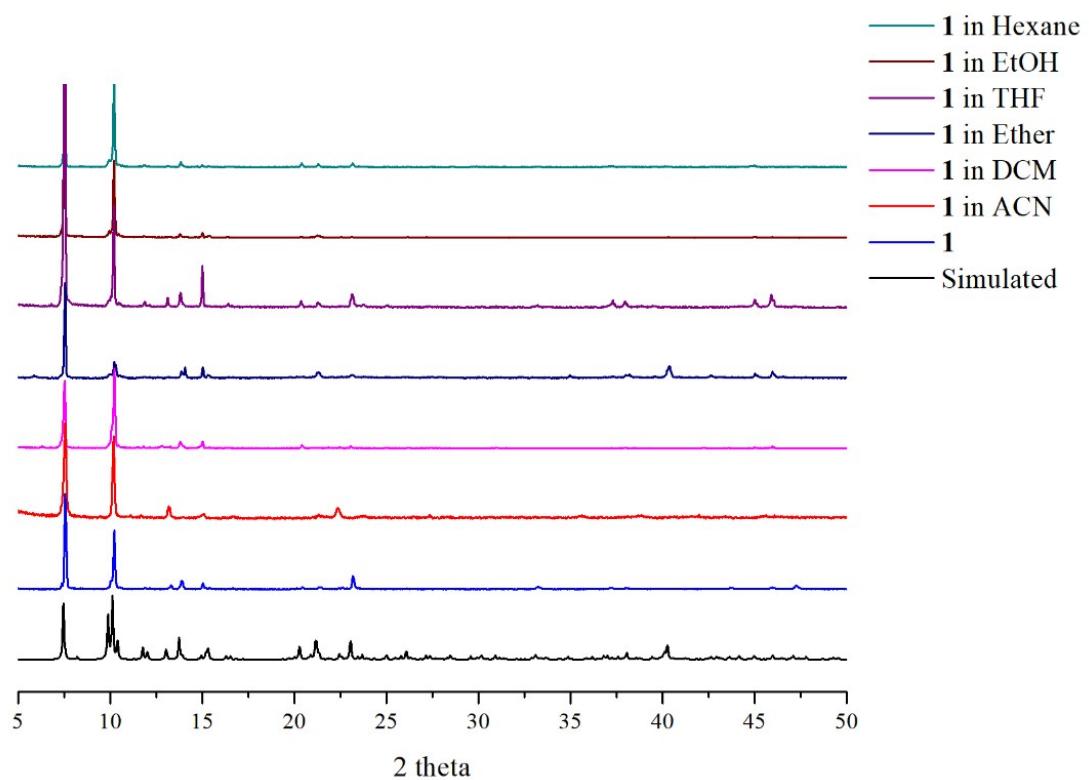


Fig. S5. (a) Complex **1** (b) **1** immersed in methanol (c) **1** immersed in methanol then back in DMF (d) **1** immersed in water (e) **1** immersed in water then back in DMF under 365 nm UV light.

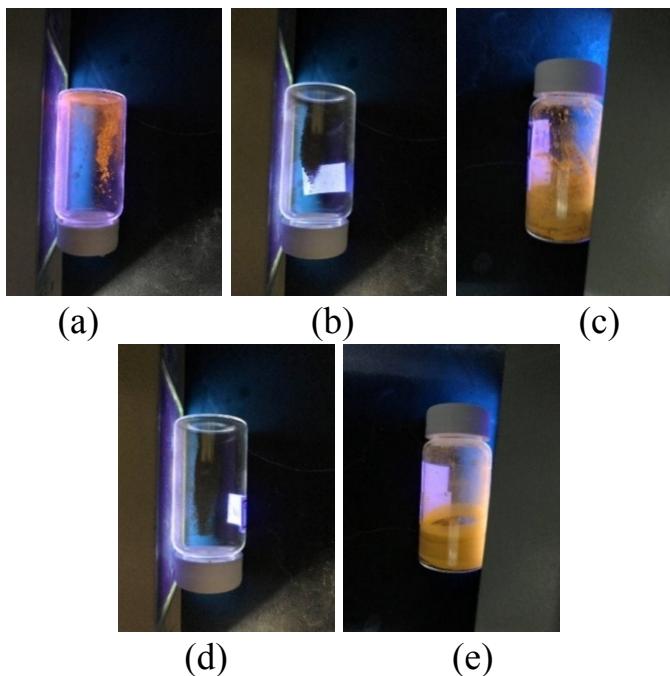


Fig. S6. PXRD patterns for **2** after immersed in various solvents.

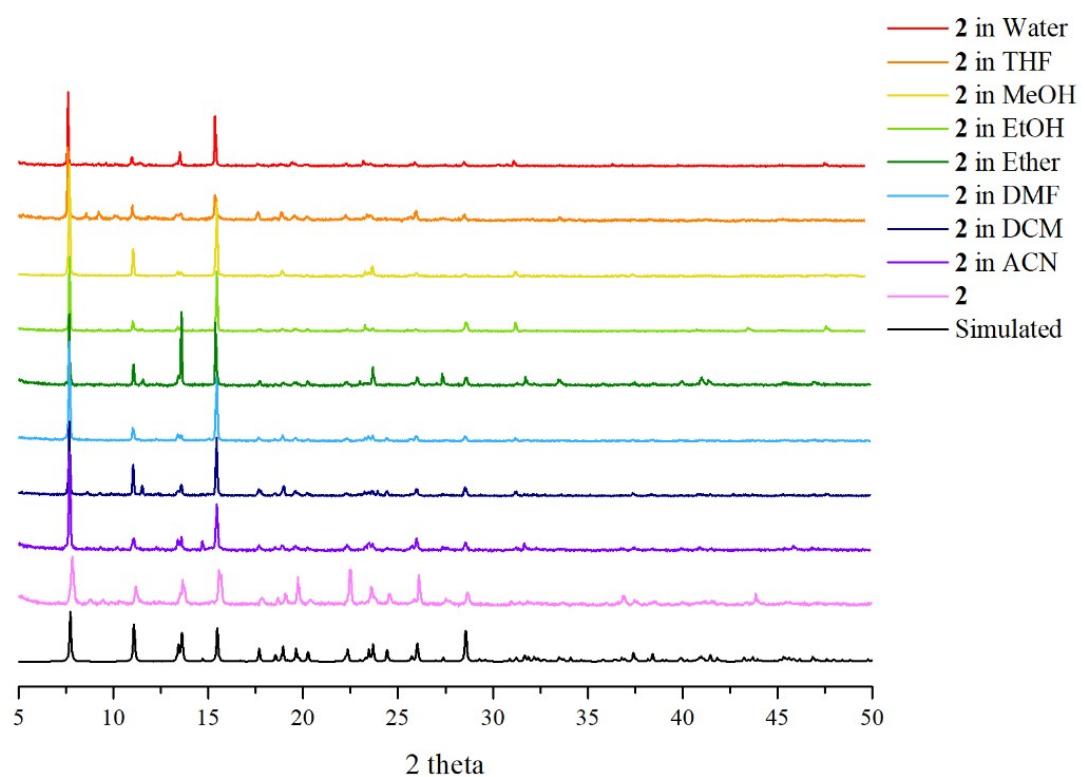


Fig. S7. TG analysis for (a) complex **1** and its hydrated product and (b) **2**.

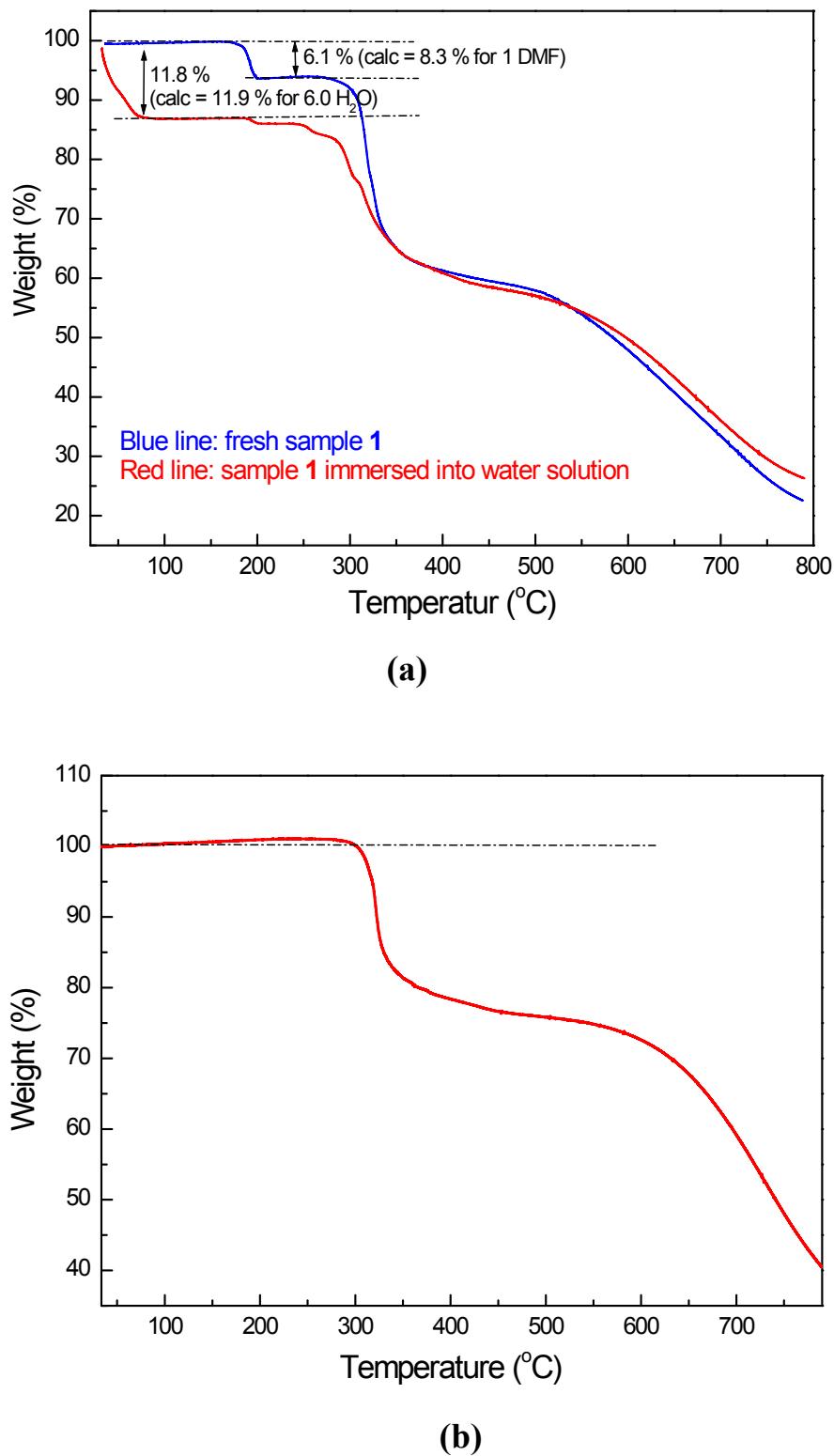


Fig. S8. TG measurements of cyclic de-/re-hydration process for **1**.

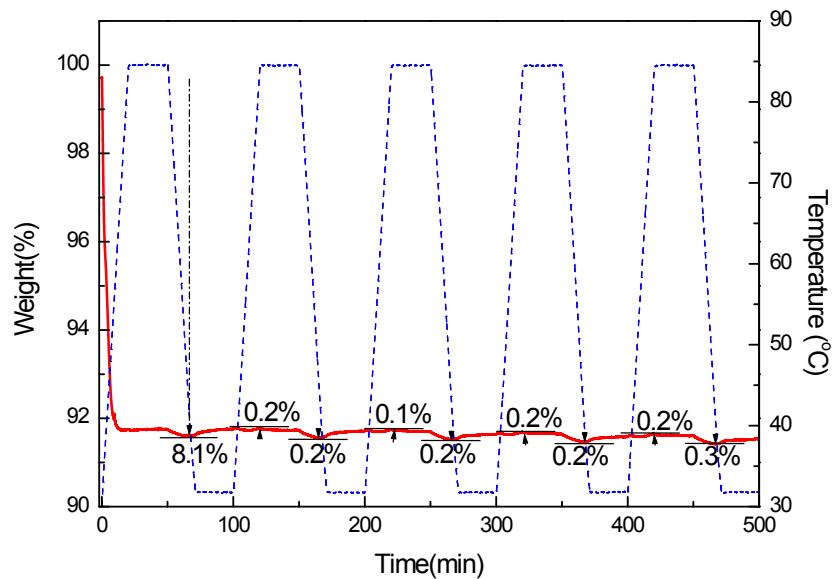


Fig. S9. Water-vapor adsorption isotherm for **1**.

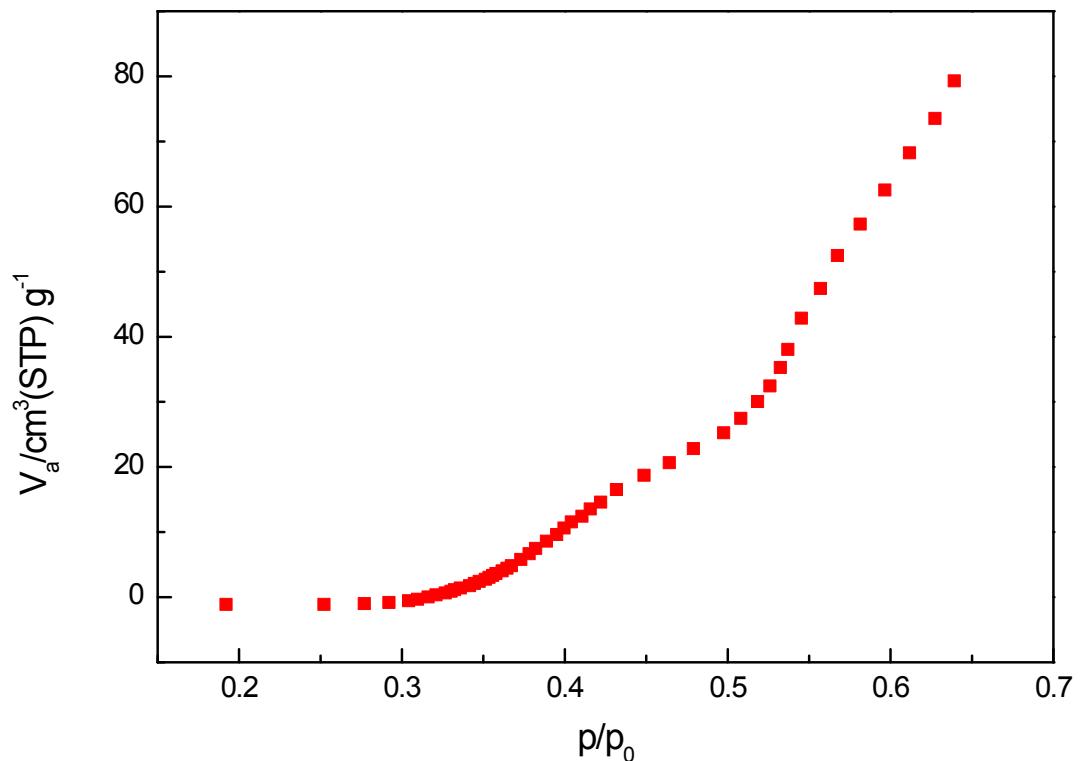
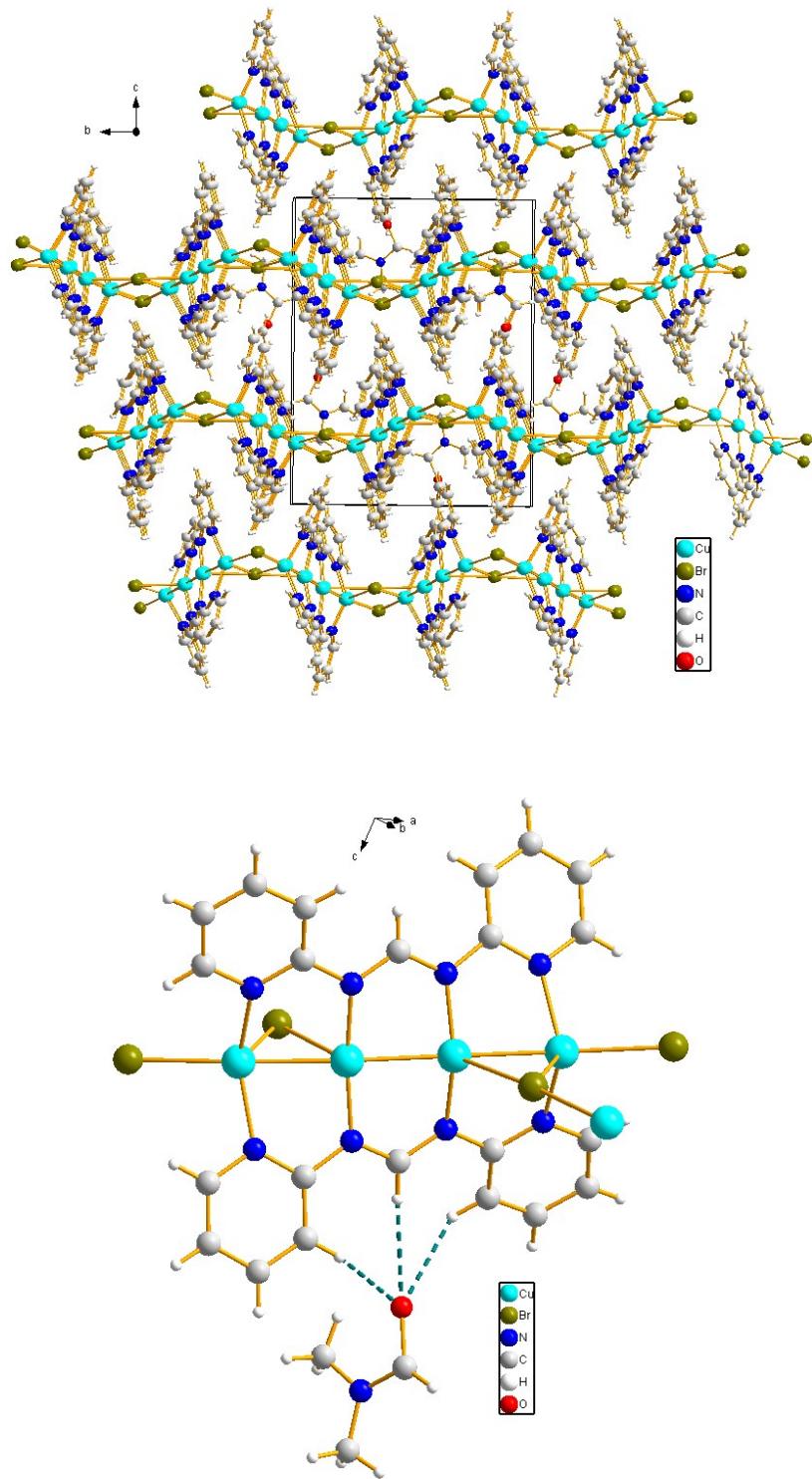


Fig. S10. Packing diagram for **1**.



EDS data for complex 1

Spectrum processing : No peaks omitted

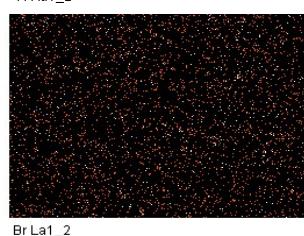
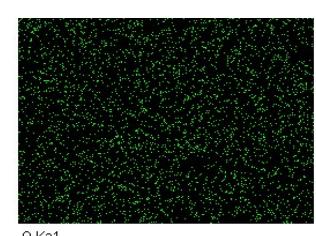
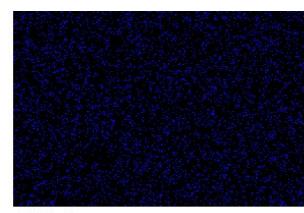
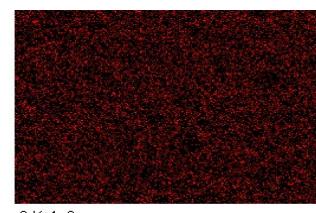
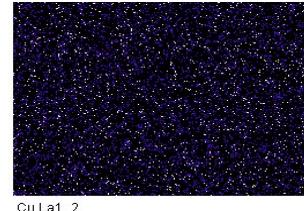
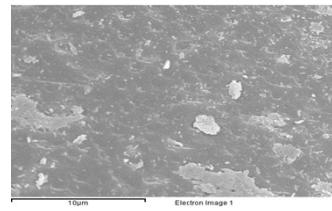
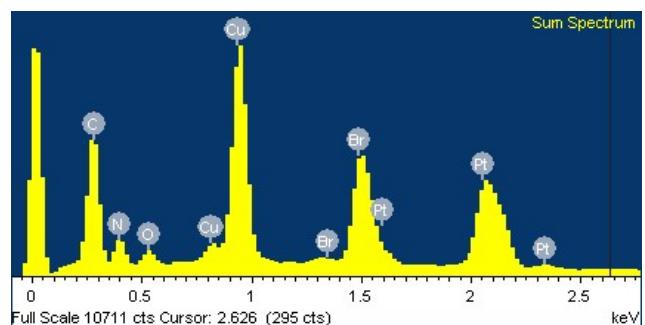
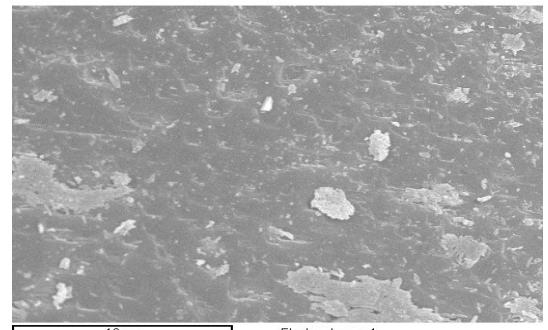
Processing option : All elements analyzed (Normalised)

Number of iterations = 3

Standard :

| | | |
|----|-------------------|---------------------|
| C | CaCO ₃ | 1-Jun-1999 12:00 AM |
| N | Not defined | 1-Jun-1999 12:00 AM |
| O | SiO ₂ | 1-Jun-1999 12:00 AM |
| Cu | Cu | 1-Jun-1999 12:00 AM |
| Br | KBr | 1-Jun-1999 12:00 AM |
| Pt | Pt | 1-Jun-1999 12:00 AM |

| Element | Weight% | Atomic% |
|---------|---------|---------|
| C K | 26.08 | 58.75 |
| N K | 9.68 | 18.70 |
| O K | 2.28 | 3.85 |
| Cu L | 26.06 | 11.10 |
| Br L | 13.10 | 4.44 |
| Pt M | 22.80 | 3.16 |
| Totals | 100.00 | |



EDS data for complex 2.

Spectrum processing : No peaks omitted

Processing option : All elements analyzed (Normalised)

Number of iterations = 3

Standard :

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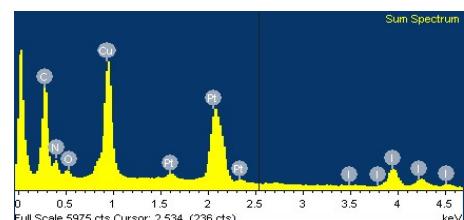
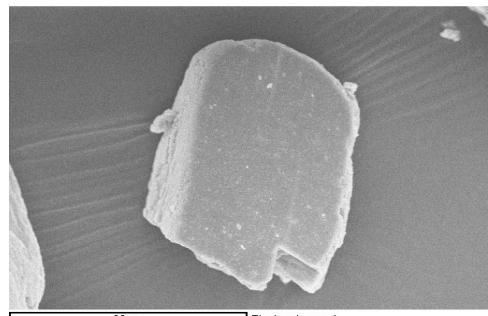
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O SiO₂ 1-Jun-1999 12:00 AM

Cu Cu 1-Jun-1999 12:00 AM

I Not defined 1-Jun-1999 12:00 AM

Pt Pt 1-Jun-1999 12:00 AM



| Element | Weight% | Atomic% |
|---------|---------|---------|
| C K | 22.90 | 60.54 |
| N K | 6.94 | 15.74 |
| O K | 1.40 | 2.77 |
| Cu L | 24.18 | 12.09 |
| I L | 18.30 | 4.58 |
| Pt M | 26.28 | 4.28 |
| Totals | 100.00 | |

