

Supporting Information

Synthesis and crystal structure of the europium(II) hydride oxide iodide $\text{Eu}_5\text{H}_2\text{O}_2\text{I}_4$ showing blue-green luminescence

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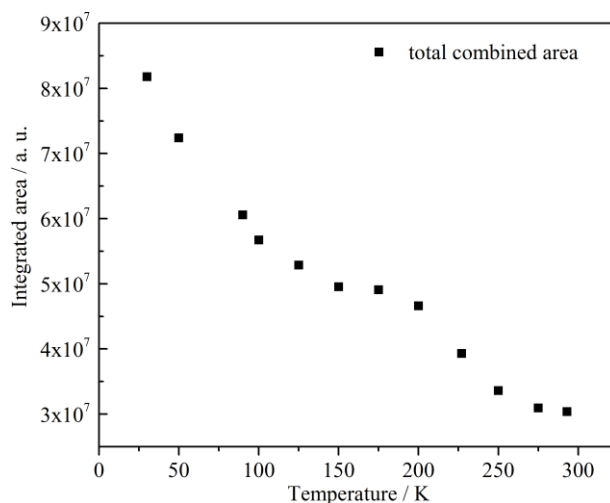


Figure S1. Temperature-dependent intensity of the emission (peak area) of $\text{Eu}_5\text{H}_2\text{O}_2\text{I}_4$. The quenching temperature is estimated to be around 110 K.

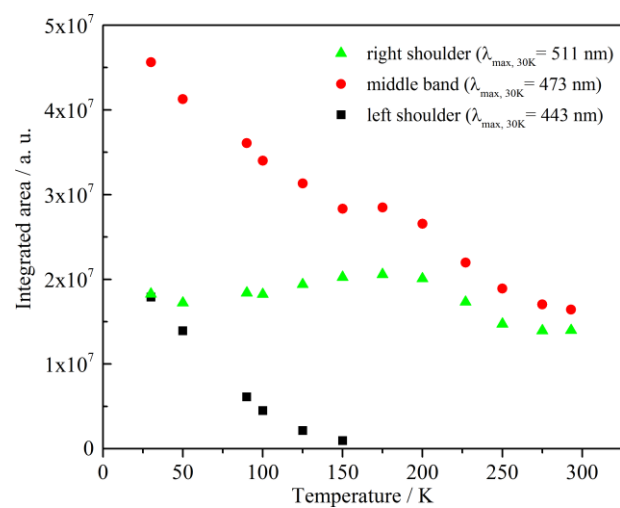


Figure S2. Integrated area of the three emission bands of $\text{Eu}_5\text{H}_2\text{O}_2\text{I}_4$ resulting after deconvolution.

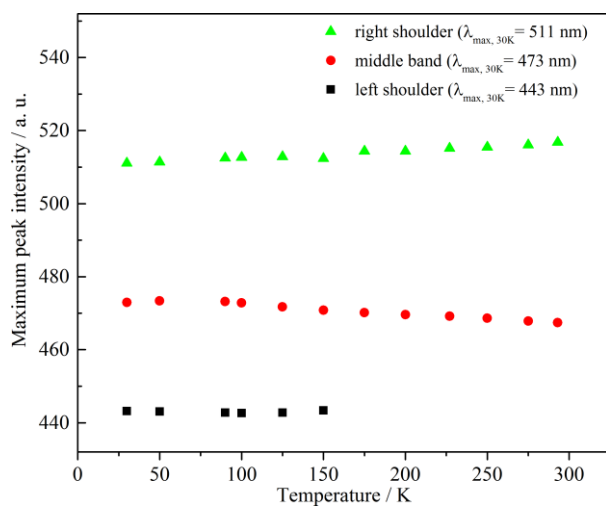


Figure S3. Change of the three individual emission positions in $\text{Eu}_5\text{H}_2\text{O}_2\text{I}_4$ with temperature.

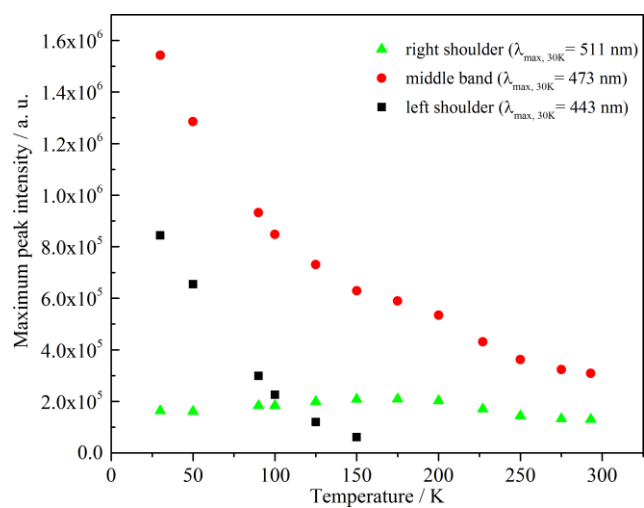


Figure S4. Maximum peak intensity of the three emission bands in $\text{Eu}_5\text{H}_2\text{O}_2\text{I}_4$ after deconvolution.

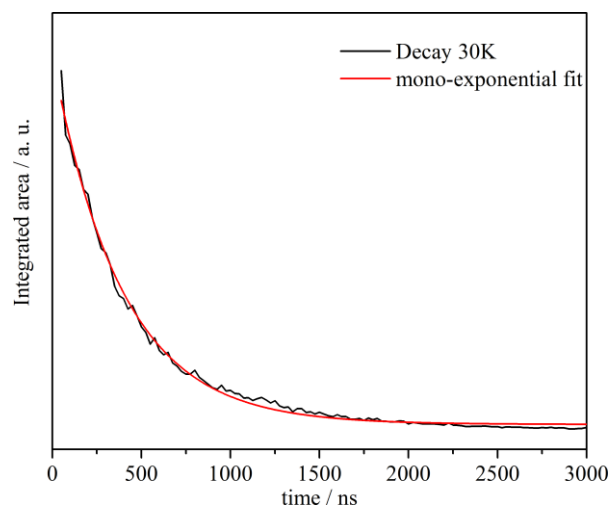


Figure S5. Decay curve of $\text{Eu}_5\text{H}_2\text{O}_2\text{I}_4$ at 30 K. Fitting of the parameters using a mono-exponential decay results in a lifetime of the excited state of 388 ± 5 ns.

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'x+1/2, -y+1/2, -z'

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'-x, -y, -z'

'x, y, -z-1/2'

'-x, y, z'

'x, -y, z-1/2'

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Refinement of F^2 against ALL reflections. The weighted R-factor wR and goodness of fit S are based on F^2 , conventional R-factors R are based on F, with F set to zero for negative F^2 . The threshold expression of $F^2 > 2\sigma(F^2)$ is used only for calculating R-factors(gt) etc. and is not relevant to the choice of reflections for refinement. R-factors based on F^2 are statistically about twice as large as those based on F, and R-factors based on ALL data will be even larger.

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O 0.021(4) 0.016(4) 0.026(4) -0.006(3) 0.000 0.000

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All esds (except the esd in the dihedral angle between two l.s. planes)
are estimated using the full covariance matrix. The cell esds are taken
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used when they are defined by crystal symmetry. An approximate (isotropic)
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'x, y, -z-1/2'

'-x, y, z'

'x, -y, z-1/2'

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Refinement of F^2 against ALL reflections. The weighted R-factor wR and goodness of fit S are based on F^2 , conventional R-factors R are based on F, with F set to zero for negative F^2 . The threshold expression of $F^2 > 2\sigma(F^2)$ is used only for calculating R-factors(gt) etc. and is not relevant to the choice of reflections for refinement. R-factors based on F^2 are statistically about twice as large as those based on F, and R-factors based on ALL data will be even larger.

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_refine_ls_weighting_details
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_atom_sites_solution_secondary  difmap
_refine_ls_extinction_method     SHELXL
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;

All esds (except the esd in the dihedral angle between two l.s. planes)
 are estimated using the full covariance matrix. The cell esds are taken
 into account individually in the estimation of esds in distances, angles
 and torsion angles; correlations between esds in cell parameters are only
 used when they are defined by crystal symmetry. An approximate (isotropic)
 treatment of cell esds is used for estimating esds involving l.s. planes.

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