Supporting Information for

Rapid Energy Transfer in Non-Porous Metal-organic Frameworks with Caged Ru(bpy)₃²⁺ Chromophores: Oxygen Trapping and Luminescence Quenching

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Figure S1. SEM images of 1.



Figure S2. SEM images of **2**.

Compound	
Empirical formula	C36 H24 N6 O12 Ru Zn2
Formula weight	964.42
Temperature (K)	100(2) K
Wavelength (Å)	1.54178
Crystal system	Cubic
Space group	P4 ₁ 32
Unit cell dimensions	a = 15.2853(5) Å
	b = 15.2853(5) Å

	c = 15.2853(5) A
	$\alpha = 90^{\circ}$
	$\beta = 90^{\circ}$
	$\gamma = 90$ °
Volume (Å ³)	3571.3(2)
Z	4
Density (calcd. g/cm ³)	1.794
Absorption coeff. (mm ⁻¹)	5.602
F(000)	1928
Crystal size (mm)	0.5×0.1 ×0.1
Crystal color & shape	red
θ range data collection	4.09 to 72.35 °
Limiting indices	-12 <u>≤</u> <i>h</i> ≤18
	-18≤ <i>k</i> ≤15
	-18≤ <i>l</i> ≤12
Reflections collected	13207
Independent reflections	1193 [$R(int) = 0.0259$]
Refinement method	Full-matrix least-square on F ²
Data/restraints/parameters	1193 / 0 / 88
Goodness-of-fit on F ²	0.795
Final R indices $[I>2\sigma(I)]^{a,b}$	R1 = 0.0220
	wR2 = 0.0799
R indices (all data)	R1 = 0.0220
	wR2 = 0.0799

 ${}^{a}R(F) = \Sigma ||F_{o}| - |F_{c}||/\Sigma |F_{o}|. {}^{b}R_{W}(F^{2}) = [\Sigma \{w(F_{o}^{2} - F_{c}^{2})^{2}\}/\Sigma \{w(F_{o}^{2})^{2}\}]^{0.5}; w^{-1} = \sigma^{2}(F_{o}^{2}) + (aP)^{2} + bP, \text{ where } P = [F_{o}^{2} + 2F_{c}^{2}]/3 \text{ and } a \text{ and } b \text{ are constants adjusted by the program.}$



Figure S3. TGA curve for 1.



Figure S4. PXRD patterns of 2 with various Os doping levels.



Figure S5. FT-IR spectrum of 1.

Table S2. Biexponential fit parameters and average lifetime for Os-doped 2.

2-xOs	A_1	τ_1	A_2	τ_2	$< \tau >$ Lifetime (ns)
0.00					1305
0.002	15.2	134	84.8	319	306
0.01	53.8	44	46.2	99.9	81
1.00					53

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1-xOs	A_1	$ au_1$	A_2	τ_2	$< \tau >$ Lifetime (ns)
0.00	20	285	80	802	760
0.002	15.6	106	84.4	271	260
0.01	26.3	31.4	73.7	85.9	79.6
1.00	13.8	19	86.2	50.3	48.5