

Supporting Information for

**Tetraplatinum Cluster Complexes Bearing Hydrophilic Anchors as
Precursors for g-Al₂O₃-Supported Platinum Nanoparticles**

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Tables of crystal data and refinement parameters for **2a**

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Table S1. Crystallographic Data of **2a**^{a,b}

complex	2a
empirical formula	C ₃₆ H ₃₂ O ₂₀ Pt ₄ ·H ₂ O·1.75(Et ₂ O)
formula weight	1710.71
temperature, K	113(1)
cryst. System	monoclinic
space group	<i>Cc</i> (#9)
<i>a</i> , Å	30.106(3)
<i>b</i> , Å	9.3586(8)
<i>c</i> , Å	23.564(2)
α , deg	-
β , deg	128.220(2)
γ , deg	-
volume, Å ³	5216.0(8)
<i>Z</i>	4
calculated density, Mg m ⁻³	2.178
2 θ_{max} , deg	54.8
limiting indices	-38 ≤ <i>h</i> ≤ 38, -12 ≤ <i>k</i> ≤ 12, -30 ≤ <i>l</i> ≤ 30
absorption coefficient, mm ⁻¹	10.792
F(000)	3206.00
crystal size, mm	0.280 x 0.210 x 0.050
goodness-of-fit on <i>F</i> ²	1.155
reflections collected / unique	39354 / 10880 [<i>R</i> (int) = 0.0686]
no. of variables	615
<i>R</i> 1 [<i>I</i> > 2 σ (<i>I</i>)]	0.0390
<i>R</i> 1, <i>wR</i> 2 (all data)	0.0609, 0.1041
residual electron density, e Å ⁻³	1.770 (max), -1.740 (min)

^a *R*1 = (Σ||*F*_o| - |*F*_c||) / Σ|*F*_o|. ^b *wR*2 = [Σ*w*(*F*_o²-*F*_c²)² / Σ(*wF*_o⁴)]^{1/2}; The function minimized: *w*(*F*_o²- *F*_c²)².