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SUPPLEMENTARY MATERIAL TO
**Dioxidovanadium(V) complexes with pyridoxal aminoguanidine
derivative: synthesis and spectral and structural
characterization**

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TABLE S-I. Hydrogen-bonding geometry parameters in compound **1**

Bond (<i>D</i> – <i>H</i> ··· <i>A</i>)	Distances, Å		Angle, °
	<i>H</i> ··· <i>A</i>	<i>D</i> ··· <i>A</i>	
O2–H2···O4 ^a	1.95	2.851(3)	163.7
N1–H1···O2 ^b	2.34	3.105(3)	144.9
N4–H4A···O4 ^c	2.25	3.136(3)	168.3
N4–H4B···O5 ^d	1.99	2.870(3)	170.9
O5–H5B···N2 ^e	1.99	2.861(3)	170.1
O5–H5A···N5	1.85	2.772(3)	168.4
N6–H6A···O3 ^f	2.15	2.862(3)	162
N6–H6B···O3	1.94	2.850(3)	161.2
N6–H6C···N2 ^b	2.02	3.003(3)	175.7
N6–H6D···O5 ^g	2.09	2.835(3)	144.3

^a_{x-1, y, z}; ^b_{-x+1, -y, -z+2}; ^c_{-x+2, -y, -z+2}; ^d_{x+1, y-1, z+1}; ^e_{-x+1, -y, -z+1}; ^f_{-x+1, -y+1, -z+2}; ^g_{-x, -y+1, -z+1}, are the equivalent positions

TABLE S-II. Hydrogen-bonding geometry parameters in compound **3**

Bond (<i>D</i> – <i>H</i> ··· <i>A</i>)	Distances, Å		Angle, °
	<i>H</i> ··· <i>A</i>	<i>D</i> ··· <i>A</i>	
O2–H2···O4 ^a	2.03	2.885(3)	164.4
N1–H1···O2 ^b	2.33	3.071(3)	149.4
N4–H4B···O5 ^c	1.98	2.813(3)	165.2
N4–H4A···O4 ^d	2.31	3.121(3)	175.6
O5–H5B···N2 ^e	1.88	2.776(3)	172.6
O5–H5A···N5	1.89	2.782(3)	167

^a_{x-1, y, z}; ^b_{-x+1, -y, -z+2}; ^c_{x+1, y-1, z+1}; ^d_{-x+2, -y, -z+2}; ^e_{-x+1, -y, -z+1}, are the equivalent positions

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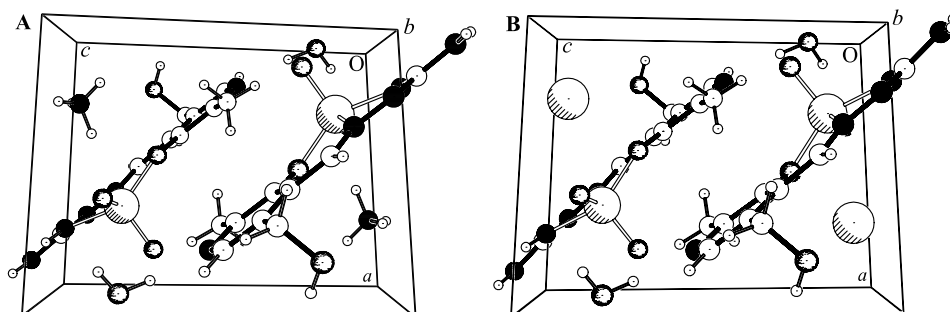


Fig. S-1. Unit cell packing for the complexes **1** (A) and **3** (B).