

Electronic Supplementary Information (ESI)

Table S1 Crystals Structure Data and Details of Structure Refinements for **3**, **4** and **5**.

Crystal data	3	4	5
CCDC deposit No.	CCDC-728424	CCDC-728423	CCDC-728425
Empirical fomular	C ₈₄ H ₈₈ N ₂ O ₄₆ S ₁₀	C ₄₇ H ₅₆ N ₄ O ₂₆ S ₅	C ₄₂ H ₄₄ N ₁ O ₂₃ S ₅
Fomular weight	2182.16	1253.26	1091.08
Crystal size/mm	0.30 × 0.26 × 0.20	0.24 × 0.20 × 0.16	0.24 × 0.22 × 0.16
Crystal system	Monoclinic	Monoclinic	Monoclinic
Space group	C2/c	P2 ₁ /n	C2/m
<i>a</i> /Å	31.666(5)	11.4886(19)	31.878(4)
<i>b</i> /Å	19.154(3)	19.653(4)	19.919(2)
<i>c</i> /Å	22.704(4)	29.851(5)	11.1685(13)
α /°	90	90	90
β /°	111.471(3)	99.170(3)	96.582(2)
γ /°	90	90	90
Volume/Å ³	12816(4)	6654(2)	7044.9(14)
Z	4	4	4
<i>D</i> _{calc} /g cm ⁻³	1.131	1.251	1.029
<i>F</i> (000)	4536	2616	2268
μ (Mo K α)/mm ⁻¹	0.246	0.250	0.224
Temperature/K	294(2)	113(2)	294(2)
<i>R</i> _{int}	0.0624	0.0535	0.0422
Range of <i>h</i> , <i>k</i> , <i>l</i>	-38/39, -23/23, -28/17	-9/13, -23/22, -35/35	-37/25, -23/23, -13/13
θ _{min/max} /°	1.27/26.51	3.13/25.00	1.29/25.01
Reflections			
collected/unique	34434/13180	37529/11636	17640/6308
Data/restraints/parameters	13180/164/710	11636/114/798	6308/445/398
Goodness of fit on <i>F</i> ²	0.854	1.022	1.347
Final <i>R</i> indices [<i>I</i> > 2 σ (<i>I</i>)]	<i>R</i> ₁ = 0.0810 <i>wR</i> ₂ = 0.2109	<i>R</i> ₁ = 0.1002 <i>wR</i> ₂ = 0.2710	<i>R</i> ₁ = 0.1508 <i>wR</i> ₂ = 0.3620
<i>R</i> indices (all data)	<i>R</i> ₁ = 0.1942 <i>wR</i> ₂ = 0.2588	<i>R</i> ₁ = 0.1207 <i>wR</i> ₂ = 0.2875	<i>R</i> ₁ = 0.2369 <i>wR</i> ₂ = 0.4337