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Table 2. Atomic coordinates ($\times 10^4$) and equivalent isotropic displacement parameters ($\text{\AA}^2 \times 10^3$) for 7. $U(\text{eq})$ is defined as one third of the trace of the orthogonalized U_{ij} tensor.

	x	y	z	U(eq)
Molecule I				
Mo(1)	2597	1528(1)	8752	33(1)
Mo(2)	2667(1)	-1072(1)	9124(1)	35(1)
Si(1)	3184(2)	-441(3)	6868(2)	55(1)
N(1)	3553(5)	357(7)	9238(4)	35(1)
N(2)	1702(5)	107(8)	8704(4)	38(2)
O(1)	2347(5)	2562(7)	9576(4)	55(2)
O(2)	2447(5)	-1445(8)	10151(4)	55(2)
C(11)	2973(7)	1271(9)	7269(5)	39(2)
C(12)	2132(7)	1947(10)	7258(6)	45(2)
C(13)	2275(9)	3211(11)	7640(8)	65(3)
C(14)	3174(9)	3343(10)	7922(7)	59(3)
C(15)	3612(6)	2154(10)	7683(6)	46(2)
C(21)	3067(7)	-1787(9)	7692(6)	47(2)
C(22)	2255(8)	-2447(10)	7863(7)	57(3)
C(23)	2422(11)	-3309(11)	8560(10)	74(4)
C(24)	3320(10)	-3256(10)	8824(8)	64(3)
C(25)	3710(7)	-2319(10)	8296(6)	48(2)
C(51)	4374(10)	-438(15)	6564(10)	84(4)
C(52)	2373(12)	-800(17)	5944(12)	95(5)
C(71)	4475(6)	502(9)	9695(6)	44(2)
C(72)	4518(9)	1976(13)	10092(9)	70(3)
C(73)	5216(7)	370(11)	9122(7)	53(2)
C(74)	4571(9)	-514(13)	10411(8)	69(3)
C(81)	709(7)	51(11)	8682(7)	53(2)
C(82)	234(19)	-112(27)	7765(16)	166(9)
C(83)	382(18)	1298(22)	9121(17)	137(8)
C(84)	432(13)	-1068(16)	9221(12)	98(5)
Molecule II				
Mo(3)	6586(1)	-5919(1)	8075(1)	33(1)
Mo(4)	6774(1)	-3339(1)	7719(1)	32(1)
Si(2)	8821(2)	-5231(3)	7284(2)	55(1)
N(3)	6322(5)	-4854(7)	7044(4)	35(1)
N(4)	6983(5)	-4385(8)	8787(5)	39(2)
O(3)	5564(5)	-6334(7)	8409(5)	53(2)
O(4)	5851(5)	-2384(7)	7869(5)	52(2)
C(31)	8105(6)	-6584(9)	7700(6)	40(2)
C(32)	7471(7)	-7402(9)	7209(6)	46(2)
C(33)	7021(8)	-8219(10)	7784(8)	54(3)
C(34)	7374(8)	-7896(12)	8631(7)	62(3)
C(35)	8028(7)	-6926(11)	8580(6)	51(2)
C(41)	8296(5)	-3537(9)	7278(6)	37(2)
C(42)	7757(7)	-2877(13)	6597(7)	60(3)
C(43)	7485(8)	-1600(12)	6904(10)	69(4)
C(44)	7849(7)	-1492(11)	7770(10)	63(3)
C(45)	8327(6)	-2658(10)	7978(7)	48(2)
C(61)	9050(12)	-5703(17)	6198(11)	95(4)
C(62)	9876(12)	-5110(18)	8034(12)	95(4)
C(91)	5758(7)	-5076(11)	6223(7)	54(2)
C(92)	5216(11)	-6357(15)	6235(11)	84(4)
C(93)	6368(10)	-5290(13)	5431(9)	73(3)
C(94)	5196(12)	-3922(15)	5975(11)	89(4)
C(101)	7015(8)	-4068(10)	9706(7)	52(2)
C(102)	6140(16)	-3488(23)	9911(17)	135(7)
C(103)	7680(17)	-3133(24)	9979(18)	160(9)
C(104)	7155(17)	-5310(20)	10281(16)	123(7)

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Table 5. Hydrogen coordinates ($\times 10^4$) and isotropic displacement parameters ($\text{\AA}^2 \times 10^3$) for 7.

	x	y	z	U(eq)
H(12A)	1582(7)	1643(10)	6959(6)	80
H(13A)	1831(9)	3910(11)	7654(8)	80
H(14A)	3458(9)	4153(10)	8149(7)	80
H(15A)	4248(6)	2016(10)	7738(6)	80
H(22A)	1700(8)	-2380(10)	7511(7)	80
H(23A)	2006(11)	-3954(11)	8756(10)	80
H(24A)	3634(10)	-3837(10)	9242(8)	80
H(25A)	4339(7)	-2136(10)	8305(6)	80
H(32A)	7411(7)	-7469(9)	6590(6)	80
H(33A)	6621(8)	-8953(10)	7625(8)	80
H(34A)	7254(8)	-8380(12)	9145(7)	80
H(35A)	8416(7)	-6607(11)	9061(6)	80
H(42A)	7669(7)	-3183(13)	6009(7)	80
H(43A)	7187(8)	-901(12)	6560(10)	80
H(44A)	7827(7)	-691(11)	8118(10)	80
H(45A)	8684(6)	-2784(10)	8514(7)	80
H(51A)	4435(10)	255(15)	6140(10)	80
H(51B)	4787(10)	-271(15)	7058(10)	80
H(51C)	4499(10)	-1303(15)	6321(10)	80
H(52A)	1781(12)	-787(17)	6138(12)	80
H(52B)	2416(12)	-127(17)	5503(12)	80
H(52C)	2491(12)	-1676(17)	5712(12)	80
H(61A)	9421(12)	-5024(17)	5967(11)	80
H(61B)	9344(12)	-6565(17)	6188(11)	80
H(61C)	8488(12)	-5747(17)	5854(11)	80
H(62A)	10251(12)	-4416(18)	7823(12)	80
H(62B)	9736(12)	-4888(18)	8609(12)	80
H(62C)	10186(12)	-5960(18)	8040(12)	80
H(72A)	5094(9)	2115(13)	10399(9)	80
H(72B)	4424(9)	2644(13)	9645(9)	80
H(72C)	4060(9)	2056(13)	10486(9)	80
H(73A)	5194(7)	-516(11)	8868(7)	80
H(73B)	5145(7)	1040(11)	8673(7)	80
H(73C)	5783(7)	496(11)	9447(7)	80
H(74A)	4548(9)	-1403(13)	10162(8)	80
H(74B)	5134(9)	-390(13)	10743(8)	80
H(74C)	4092(9)	-413(13)	10781(8)	80
H(82A)	-405(19)	-151(27)	7783(16)	80
H(82B)	388(19)	654(27)	7428(16)	80
H(82C)	439(19)	-926(27)	7508(16)	80
H(83A)	-259(18)	1296(22)	9113(17)	80
H(83B)	637(18)	1358(22)	9707(17)	80
H(83C)	571(18)	2060(22)	8799(17)	80
H(84A)	-210(13)	-1104(16)	9197(12)	80
H(84B)	657(13)	-1898(16)	9002(12)	80
H(84C)	669(13)	-943(16)	9809(12)	80
H(92A)	4822(11)	-6329(15)	6692(11)	80
H(92B)	4871(11)	-6481(15)	5691(11)	80
H(92C)	5628(11)	-7096(15)	6331(11)	80
H(93A)	6727(10)	-4498(13)	5367(9)	80
H(93B)	6751(10)	-6058(13)	5547(9)	80
H(93C)	5995(10)	-5443(13)	4908(9)	80
H(94A)	4809(12)	-3735(15)	6422(11)	80
H(94B)	5581(12)	-3159(15)	5907(11)	80
H(94C)	4842(12)	-4088(15)	5441(11)	80
H(10A)	6140(16)	-3270(23)	10513(17)	80
H(10B)	6048(16)	-2679(23)	9573(17)	80
H(10C)	5667(16)	-4119(23)	9754(17)	80
H(10D)	7673(17)	-2966(24)	10588(18)	80
H(10E)	8254(17)	-3484(24)	9860(18)	80
H(10F)	7568(17)	-2303(24)	9668(18)	80
H(10G)	7173(17)	-5063(20)	10881(16)	80
H(10H)	6685(17)	-5956(20)	10152(16)	80
H(10I)	7718(17)	-5701(20)	10162(16)	80

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Table 4. Anisotropic displacement parameters ($\text{\AA}^2 \times 10^3$) for 7.
 The anisotropic displacement factor exponent takes the form:
 $-2\pi^2 [h^2 a^{*2} U_{11} + \dots + 2 h k a^* b^* U_{12}]$

	U11	U22	U33	U23	U13	U12
Molecule I						
Mo(1)	37(1)	35(1)	28(1)	-5(1)	8(1)	2(1)
Mo(2)	38(1)	38(1)	30(1)	3(1)	7(1)	-4(1)
Si(1)	98(2)	44(1)	24(1)	-2(1)	12(1)	16(1)
N(1)	38(4)	38(4)	29(3)	-1(3)	10(3)	-2(3)
N(2)	36(4)	53(4)	26(3)	0(3)	5(3)	-5(3)
O(1)	66(4)	60(4)	40(3)	-18(3)	14(3)	11(4)
O(2)	58(4)	68(4)	40(3)	13(3)	15(3)	-10(4)
C(11)	60(5)	44(4)	16(3)	6(3)	13(3)	5(4)
C(12)	49(5)	56(5)	32(4)	14(4)	6(4)	5(4)
C(13)	85(8)	48(6)	64(7)	18(5)	22(6)	26(5)
C(14)	88(9)	39(5)	54(6)	2(4)	19(6)	-12(5)
C(15)	44(5)	56(5)	39(4)	9(4)	13(4)	-11(4)
C(21)	58(6)	44(5)	39(5)	-14(4)	12(4)	9(4)
C(22)	69(7)	43(5)	58(6)	-16(4)	-4(5)	-11(5)
C(23)	101(12)	37(5)	87(10)	-3(5)	21(8)	-28(6)
C(24)	93(9)	37(5)	64(7)	9(5)	19(6)	13(5)
C(25)	51(5)	48(5)	46(5)	-5(4)	15(4)	14(4)
Molecule II						
Mo(3)	34(1)	37(1)	31(1)	2(1)	7(1)	0(1)
Mo(4)	29(1)	33(1)	35(1)	-2(1)	6(1)	4(1)
Si(2)	38(1)	48(1)	82(2)	-11(1)	24(1)	3(1)
N(3)	32(3)	38(3)	34(3)	-1(3)	7(3)	2(3)
N(4)	36(4)	50(4)	32(3)	0(3)	8(3)	8(3)
O(3)	41(4)	63(4)	58(4)	15(3)	18(3)	-7(3)
O(4)	46(3)	52(4)	60(4)	-6(3)	11(3)	17(3)
C(31)	34(4)	42(4)	46(5)	-8(4)	11(4)	11(3)
C(32)	61(6)	33(4)	46(5)	-9(4)	19(4)	10(4)
C(33)	59(7)	39(5)	64(6)	-1(4)	15(5)	2(4)
C(34)	72(8)	61(6)	53(6)	26(5)	14(5)	14(6)
C(35)	51(5)	56(5)	45(5)	-3(4)	-3(4)	18(5)
C(41)	24(4)	48(4)	42(4)	1(4)	12(3)	-1(3)
C(42)	48(5)	84(8)	50(6)	20(5)	23(4)	-17(5)
C(43)	51(6)	55(6)	105(10)	35(6)	22(7)	-2(5)
C(45)	34(4)	51(5)	57(6)	-7(4)	0(4)	-15(4)
C(44)	41(5)	43(5)	107(10)	-9(6)	12(6)	-9(4)