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Table. Non-hydrogen positional and isotropic displacement parameters

Atom	*x*	*y*	*z*	U_{eq} Å^2
Ag(1)	0.2227(1)	0.7929(1)	0.49826(3)	0.0557(6)
Ag(2)	0.2239(1)	X+1	1/2	0.0623(7)
Ag(3)	-0.3274(1)	X+2	1/2	0.0537(6)
O	-0.025(1)	1.247(1)	0.4523(2)	0.038(4)
C(1)	-0.002(2)	1.136(2)	0.4341(3)	0.048(7)
C(2)	0.041(2)	1.006(2)	0.4495(3)	0.047(7)
S(3)	0.2059(5)	1.0098(5)	0.46828(9)	0.042(2)
C(4)	0.339(2)	1.037(2)	0.4403(4)	0.063(9)
C(5)	0.362(2)	0.919(2)	0.4233(5)	0.07(1)
N(6)	0.499(2)	0.936(2)	0.4065(3)	0.073(7)
C(1')	-0.050(2)	1.368(2)	0.4367(4)	0.056(9)
C(2')	-0.098(2)	1.485(2)	0.4555(4)	0.048(7)
S(3')	-0.2585(5)	1.4493(4)	0.47523(9)	0.035(2)
C(4')	-0.380(2)	1.400(2)	0.4478(3)	0.035(7)
C(5')	-0.419(2)	1.519(2)	0.4284(3)	0.045(7)
N(6')	-0.530(1)	1.472(1)	0.4083(3)	0.049(6)
N(1)	-0.033(2)	0.679(2)	0.5261(3)	0.052(7)
O(11)	-0.036(1)	0.739(1)	0.5030(3)	0.073(6)
O(12)	0.071(1)	0.681(1)	0.5416(3)	0.069(6)
O(13)	-0.140(1)	0.615(1)	0.5349(2)	0.059(5)
N(2)	0.496(2)	1.261(1)	0.5301(3)	0.043(6)
O(21)	0.389(1)	1.287(1)	0.5429(3)	0.062(5)
O(22)	0.612(1)	1.277(2)	0.5424(3)	0.073(6)
O(23)	0.495(1)	1.218(1)	0.5049(3)	0.066(5)
N(3)	-0.133(3)	0.779(3)	0.3948(5)	0.11(1)
O(31)	-0.192(2)	0.832(2)	0.4136(4)	0.15(1)
O(32)	-0.033(2)	0.807(2)	0.3808(4)	0.126(9)
O(33)	-0.162(3)	0.659(3)	0.3917(5)	0.24(2)
N(4)	0.211(3)	0.676(3)	0.6235(4)	0.09(1)
O(41)	0.329(2)	0.725(2)	0.6246(4)	0.116(9)
O(42)	0.207(2)	0.560(2)	0.6098(3)	0.088(7)
O(43)	0.105(2)	0.719(2)	0.6324(4)	0.16(1)