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Chlorido{4-ethyl-1-[1-(pyrazin-2-yl)ethylidene]thiosemicabazidato-κS}bis-(triphenylphosphane-κP)silver(I)

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Key indicators: single-crystal X-ray study; T = 90 K; mean σ (C–C) = 0.005 Å; R factor = 0.042; wR factor = 0.103; data-to-parameter ratio = 27.4.

The title compound, $[Ag(C_9H_{13}N_5S)Cl(C_{18}H_{15}P)_2]$, crystallizes with four independent molecules in the asymmetric unit, in each of which the Ag atom is in a distorted tetrahedral coordination, defined by the chloride ligand, the S atom of the neutral ligand and two P atoms derived from the triphenyl phosphine ligands. The thiosemicarbazone acts as a monodentate ligand through its thione S atom. An intramolecular $N-H\cdots$ Cl hydrogen bond occurs in two of the independent molecules. In the crystal, the molecules are assembled through $N-H\cdots$ Cl hydrogen bonds, forming chains along [101].

Related literature

For general background to thiosemicarbazones, see: Akinchan *et al.* (2002); Ali & Livingstone (1974); Bermejo *et al.* (2003); Blanz & French (1968); Campbell (1975); Casas *et al.* (2000); Grecu & Neamtu (1967); Hossain *et al.* (2002); Huheey *et al.* (1993); Lobana *et al.* (1998, 2008); Pellerito & Negy (2002); Raper (1985); Venkatraman *et al.* (2009); Zhou *et al.* (2008).



Experimental

Crystal data

 $\begin{bmatrix} Ag(C_9H_{13}N_5S)Cl(C_{18}H_{15}P)_2 \end{bmatrix} \\ M_r = 891.17 \\ Monoclinic, P2_1 \\ a = 12.0640 (5) \text{ Å} \\ b = 31.810 (2) \text{ Å} \\ c = 21.9207 (15) \text{ Å} \\ \beta = 98.029 (5)^{\circ} \end{bmatrix}$

Data collection

Nonius KappaCCD diffractometer with Oxford Cryostream system Absorption correction: multi-scan (DENZO and SCALEPACK; Otwinowski & Minor, 1997) $T_{min} = 0.831, T_{max} = 0.870$

Refinement

 $R[F^2 > 2\sigma(F^2)] = 0.042$ $wR(F^2) = 0.103$ S = 1.0354942 reflections 2007 parameters 9 restraints $V = 8329.7 (9) \text{ Å}^{3}$ Z = 8 Mo K\alpha radiation $\mu = 0.71 \text{ mm}^{-1}$ T = 90 K 0.27 \times 0.22 \times 0.20 mm

138032 measured reflections 54942 independent reflections 44399 reflections with $I > 2\sigma(I)$ $R_{\text{int}} = 0.055$

H atoms treated by a mixture of
independent and constrained
refinement
$\Delta \rho_{\rm max} = 1.03 \text{ e } \text{\AA}^{-3}$
$\Delta \rho_{\rm min} = -0.85 \text{ e } \text{\AA}^{-3}$
Absolute structure: Flack (1983),
25,519 Friedel pairs
Flack parameter: 0.382 (11)

Table 1 Hydrogen-bond geometry (Å, °).

$D - H \cdot \cdot \cdot A$	D-H	$H \cdot \cdot \cdot A$	$D \cdots A$	$D - \mathbf{H} \cdot \cdot \cdot A$
$N1 - H1N \cdots Cl1 N3 - H3N \cdots Cl3 N6 - H6N \cdots Cl2 N8 - H8N \cdots Cl4 N11 - H11N \cdots Cl3 N11 - H11N - Cl3$	0.88 (2)	2.54 (2)	3.381 (3)	163 (3)
	0.87 (2)	2.68 (3)	3.406 (3)	142 (3)
	0.88 (2)	2.53 (2)	3.370 (3)	162 (3)
	0.86 (2)	2.68 (3)	3.411 (3)	143 (3)
	0.88 (2)	2.52 (2)	3.363 (3)	161 (3)
N13−H13 <i>N</i> ···Cl2	0.88 (2)	2.65 (2)	3.423 (3)	147 (3)
N16−H16 <i>N</i> ···Cl4	0.86 (2)	2.50 (2)	3.338 (3)	164 (3)
N18−H18 <i>N</i> ···Cl1 ⁱ	0.86 (2)	2.70 (3)	3.425 (3)	143 (3)

Symmetry code: (i) x - 1, y, z - 1.

Data collection: *COLLECT* (Nonius, 2000); cell refinement: *SCALEPACK* (Otwinowski & Minor, 1997); data reduction: *DENZO* (Otwinowski & Minor, 1997) and *SCALEPACK*; program(s) used to solve structure: *SIR97* (Altomare *et al.*, 1999); program(s) used to refine structure: *SHELXH* (Sheldrick, 2008); molecular graphics: *ORTEP-3 for Windows* (Farrugia, 2012); software used to prepare material for publication: *SHELXH*.

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Supplementary data and figures for this paper are available from the IUCr electronic archives (Reference: RK2386).

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Chlorido{4-ethyl-1-[1-(pyrazin-2-yl)ethylidene]thiosemicabazidato-*kS*}bis(triphenylphosphane-*kP*)silver(I)

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S1. Comment

Thiosemicarbazones are versatile ligands for several metals due to the keto–enol tautomeric form exhibited by the molecules (Ali & Livingstone, 1974; Campbell, 1975; Pellerito & Negy, 2002; Raper, 1985; Casas *et al.*, 2000; Blanz & French, 1968; Grecu & Neamtu, 1967; Lobana *et al.*, 1998; Bermejo *et al.*, 2003; Akinchan *et al.*, 2002; Zhou *et al.* 2008; Huheey *et al.*, 1993). Among the metals that complex with thiosemicarbazones, Ag(I), and Cu(I) are known to interact with thiosemicarbazones differently and the resulting compounds exhibit variable structural features depending on the conditions of their preparation. Recently, the influence of substituents at the C² carbon of thiosemicarbazones on bonding and nuclearity of silver complexes were investigated (Lobana *et al.*, 2008). Among the ligands that ligate with Ag(I) with pyrazine thiosemicarbazone received less attention.

In continuation of our study on metal complexes with thiosemicarbazones, we report herein a silver complex (Scheme 1) formed by the reaction of 2–acetylpyrazine *N*(4)–ethyl–3–thiosemicarbazone with AgCl in presence of triphenyl-phosphine as a coligand. The compound crystallizes with four independent molecules in asymmetric unit (Fig. 1). The Ag —S bond distances vary from 2.6816 (8)Å to 2.7412 (9)Å with a mean value of 2.7067 (9)Å. The Ag—S bond distance is smaller than the sum of ionic radii of Ag and S ions (2.78Å). On the other hand, the distances of Ag—P are in the range of 2.4784 (8)Å to 2.4877 (8)Å with a mean value of 2.4835 (8)Å which are smaller than that observed for Ag—S bonds. The average bond distance of Ag—P is comparable to the mononuclear silver complex of thiosemicarbazone (2.4409 (7)Å and 2.4879 (7)Å) reported by Lobana *et al.* (2008). In the asymmetric unit the four Ag—Cl bonds distances are almost similar (2.5900 (8)Å to 2.5932 (8)Å) with a mean value of 2.5921 (8)Å which is shorter than the sum of the ionic radii of silver and chloride (2.75Å). The C—S bond distances in all four units is identical to 1.704 (3)Å which is much longer than the literature value (1.6796 (9)Å) for the free ligand reported by Venkatraman *et al.* (2009). As shown in the Fig. 2, the molecules are linked with intra– and inter–molecular N—H…Cl hydrogen bonding interactions (Hossain *et al.*, 2002).

S2. Experimental

To a freshly prepared AgCl (0.143 g, 1 mmol) suspended in acetonitrile (20 ml) and was mixed with an acetonitrile solution of 2–acetylpyrazine N(4)–ethyl–3–thiosemicarbazone (0.223 g, 1 mmol,) (Venkatraman *et al.*, 2009) and the resulting mixture was stirred overnight. To the white solid formed in acetonitrile, solid triphenylphosphine (0.530 g, 2 mmol) was added in two equal aliquots to obtain a clear solution. The solution was filtered and left for crystallization at room temperature.

S3. Refinement

H atom of NH group was located in difference syntheses and refined isotropically with $U_{iso}(H) = 1.2U_{eq}(N)$. The remaining H atoms were positioned geometrically with C—H = 0.95Å, 0.99Å and 0.98Å for aromatic, methylene and methyl H atoms, respectively, and constrained to ride on their parent atoms with $U_{iso}(H) = xU_{eq}(C)$, where x = 1.5 for methyl H, and x = 1.2 for all other H atoms.



Figure 1

The molecular structure of the title compound. Displacement ellipsoids are drawn at the 50% probability level. H atoms are presented as a small spheres of arbitrary radius. Only base atoms are labeled for clarity.



Figure 2

Intra- and inter-molecular hydrogen bonding in the crystal structure of title compound viewed along a axis.

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Crystal data

$[Ag(C_9H_{13}N_5S)Cl(C_{18}H_{15}P)_2]$	F(000) = 3664
$M_r = 891.17$	$D_{\rm x} = 1.421 {\rm ~Mg~m^{-3}}$
Monoclinic, $P2_1$	Mo <i>K</i> α radiation, $\lambda = 0.71073$ Å
Hall symbol: P 2yb	Cell parameters from 28230 reflections
a = 12.0640 (5) Å	$\theta = 2.5 - 32.5^{\circ}$
b = 31.810(2) Å	$\mu = 0.71 \text{ mm}^{-1}$
c = 21.9207 (15) Å	T = 90 K
$\beta = 98.029 \ (5)^{\circ}$	Block, colourless
$V = 8329.7 (9) Å^3$	$0.27 \times 0.22 \times 0.20$ mm
Z = 8	
Data collection	
Nonius KappaCCD	138032 measured reflections
diffractometer with Oxford Cryostream system	54942 independent reflections
Radiation source: fine-focus sealed tube	44399 reflections with $I > 2\sigma(I)$
Graphite monochromator	$R_{\rm int} = 0.055$
ω - and φ -scans	$\theta_{\rm max} = 32.6^\circ, \ \theta_{\rm min} = 2.5^\circ$
Absorption correction: multi-scan	$h = -18 \rightarrow 18$
(DENZO and SCALEPACK; Otwinowski &	$k = -46 \rightarrow 48$
Minor, 1997)	$l = -32 \rightarrow 32$

 $T_{\rm min} = 0.831, T_{\rm max} = 0.870$

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Refinement

Refinement on F^2	Hydrogen site location: inferred from
Least-squares matrix: full	neighbouring sites
$R[F^2 > 2\sigma(F^2)] = 0.042$	H atoms treated by a mixture of independent
$wR(F^2) = 0.103$	and constrained refinement
S = 1.03	$w = 1/[\sigma^2(F_o^2) + (0.0471P)^2 + 3.4676P]$
54942 reflections	where $P = (F_o^2 + 2F_c^2)/3$
2007 parameters	$(\Delta/\sigma)_{\rm max} = 0.002$
9 restraints	$\Delta \rho_{\rm max} = 1.03 \text{ e } \text{\AA}^{-3}$
Primary atom site location: structure-invariant	$\Delta \rho_{\rm min} = -0.85 \text{ e } \text{\AA}^{-3}$
direct methods	Absolute structure: Flack (1983), 25,519 Friedel
Secondary atom site location: difference Fourier	pairs
map	Absolute structure parameter: 0.382 (11)

Special details

Geometry. All s.u.'s (except the s.u. in the dihedral angle between two l.s. planes) are estimated using the full covariance matrix. The cell s.u.'s are taken into account individually in the estimation of s.u.'s in distances, angles and torsion angles; correlations between s.u.'s in cell parameters are only used when they are defined by crystal symmetry. An approximate (isotropic) treatment of cell s.u.'s is used for estimating s.u.'s involving l.s. planes.

Refinement. Refinement of F^2 against ALL reflections. The weighted *R*-factor *wR* and goodness of fit *S* are based on F^2 , conventional *R*-factors *R* are based on *F*, with *F* set to zero for negative F^2 . The threshold expression of $F^2 > \sigma(F^2)$ is used only for calculating *R*-factors(gt) *etc.* and is not relevant to the choice of reflections for refinement. *R*-factors based on F^2 are statistically about twice as large as those based on *F*, and *R*-factors based on ALL data will be even larger.

Fractional atomic coordinates and isotropic or equivalent isotropic displacement parameters $(Å^2)$

	x	У	Ζ	$U_{ m iso}$ */ $U_{ m eq}$	
Agl	0.723632 (16)	0.688036 (7)	0.820344 (9)	0.01593 (5)	
Cl1	0.86926 (6)	0.65407 (2)	0.90391 (3)	0.01736 (14)	
N1	0.7674 (2)	0.57932 (9)	0.80150 (12)	0.0178 (5)	
H1N	0.803 (3)	0.6002 (8)	0.8216 (14)	0.021*	
N2	0.8198 (2)	0.55463 (9)	0.76375 (13)	0.0172 (5)	
N3	0.5999 (2)	0.56061 (9)	0.74491 (12)	0.0185 (5)	
H3N	0.639 (3)	0.5417 (9)	0.7287 (16)	0.022*	
N4	1.0907 (2)	0.52049 (10)	0.74080 (14)	0.0232 (6)	
N5	0.9582 (2)	0.47812 (10)	0.64508 (13)	0.0217 (6)	
P1	0.59953 (6)	0.74435 (3)	0.84917 (4)	0.01490 (15)	
P2	0.78028 (6)	0.69445 (3)	0.71589 (3)	0.01796 (15)	
S1	0.59041 (6)	0.62016 (3)	0.83052 (4)	0.01636 (14)	
C1	0.6544 (2)	0.58431 (10)	0.78917 (13)	0.0147 (5)	
C2	0.9254 (2)	0.54798 (10)	0.77890 (14)	0.0161 (6)	
C3	0.9931 (3)	0.56410 (12)	0.83686 (16)	0.0219 (7)	
H3A	0.9773	0.5940	0.8417	0.033*	
H3B	1.0729	0.5603	0.8344	0.033*	
H3C	0.9732	0.5485	0.8723	0.033*	
C4	0.9783 (2)	0.52177 (10)	0.73496 (15)	0.0162 (6)	
C5	1.1349 (3)	0.49938 (12)	0.69781 (17)	0.0276 (8)	
H5	1.2141	0.4982	0.7004	0.033*	
C6	1.0700 (3)	0.47899 (11)	0.64920 (16)	0.0237 (7)	
H6	1.1056	0.4655	0.6185	0.028*	

C7	0.9139 (3)	0.49924 (11)	0.68781 (15)	0.0182 (6)
H7	0.8349	0.4991	0.6864	0.022*
C8	0.4778 (2)	0.56112 (12)	0.72850 (16)	0.0244 (7)
H8A	0.4581	0.5491	0.6867	0.029*
H8B	0.4514	0.5906	0.7272	0.029*
C9	0.4185 (3)	0.53695 (13)	0.77261 (18)	0.0375 (9)
H9A	0.3375	0.5386	0.7597	0.056*
H9B	0.4369	0.5489	0.8140	0.056*
H9C	0.4422	0.5075	0.7731	0.056*
C10	0.6477 (2)	0.79894 (10)	0.85914 (13)	0.0163 (6)
C11	0.7392 (2)	0.81151 (9)	0.83082 (13)	0.0205 (6)
H11	0.7782	0.7916	0.8094	0.025*
C12	0.7730(2)	0.85361 (9)	0.83422 (14)	0.0231 (6)
H12	0.8353	0.8624	0.8153	0.028*
C13	0.7158 (3)	0.88233 (11)	0.86504 (14)	0.0245 (7)
H13	0.7375	0.9111	0.8661	0.029*
C14	0.6270 (3)	0.86977 (9)	0.89457 (14)	0.0259 (6)
H14	0.5892	0.8897	0.9165	0.031*
C15	0.5932 (2)	0.82786 (9)	0.89200 (13)	0.0226 (6)
H15	0.5330	0.8191	0.9127	0.027*
C16	0.5438 (2)	0.73372 (9)	0.92063 (13)	0.0181 (6)
C17	0.4305 (3)	0.73699 (9)	0.92650 (14)	0.0255 (6)
H17	0.3782	0.7452	0.8921	0.031*
C18	0.3938 (3)	0.72836 (10)	0.98251 (16)	0.0324 (7)
H18	0.3164	0.7301	0.9861	0.039*
C19	0.4705 (3)	0.71709 (10)	1.03343 (16)	0.0362 (8)
H19	0.4456	0.7119	1.0720	0.043*
C20	0.5821 (4)	0.71342 (13)	1.02790 (18)	0.0390 (9)
H20	0.6344	0.7059	1.0627	0.047*
C21	0.6185 (3)	0.72077 (12)	0.97114 (15)	0.0313 (8)
H21	0.6951	0.7169	0.9670	0.038*
C22	0.4748 (2)	0.74862 (10)	0.79088 (13)	0.0187 (6)
C23	0.4369 (3)	0.78667 (10)	0.76388 (14)	0.0283 (7)
H23	0.4757	0.8120	0.7758	0.034*
C24	0.3418 (3)	0.78747 (10)	0.71930 (16)	0.0351 (8)
H24	0.3160	0.8135	0.7013	0.042*
C25	0.2852 (3)	0.75088 (11)	0.70126 (15)	0.0313 (7)
H25	0.2206	0.7517	0.6711	0.038*
C26	0.3230 (3)	0.71324 (10)	0.72730(15)	0.0275(7)
H26	0.2838	0.6880	0.7153	0.033*
C27	0.4186 (2)	0.71185 (10)	0.77124 (14)	0.0237 (6)
H27	0.4454	0.6856	0.7878	0.028*
C28	0.9007 (3)	0.66345 (10)	0.70072 (15)	0.0184 (6)
C29	0.9094 (3)	0.64519 (11)	0.64350 (15)	0.0203 (6)
H29	0.8501	0.6484	0.6104	0.024*
C30	1.0044 (3)	0.62231 (11)	0.63466 (16)	0.0232 (7)
H30	1.0089	0.6095	0.5959	0.028*
C31	1.0919 (3)	0.61826 (12)	0.68209 (16)	0.0244 (7)
	× /	× /	· · · ·	

H31	1.1575	0.6034	0.6755	0.029*
C32	1.0845 (3)	0.63581 (13)	0.73915 (17)	0.0260 (7)
H32	1.1449	0.6329	0.7717	0.031*
C33	0.9888 (3)	0.65777 (11)	0.74910 (15)	0.0228 (6)
H33	0.9830	0.6689	0.7887	0.027*
C34	0.6669 (2)	0.68237 (10)	0.65440 (13)	0.0210 (6)
C35	0.5753 (3)	0.66037 (11)	0.66935 (14)	0.0266 (7)
H35	0.5752	0.6511	0.7105	0.032*
C36	0.4834 (3)	0.65156(12)	0.62534 (15)	0.0331 (8)
H36	0.4211	0.6368	0.6368	0.040*
C37	0.4826 (3)	0.66418 (10)	0.56546 (14)	0.0280(7)
H37	0.4201	0.6580	0.5354	0.034*
C38	0.5734(3)	0.68602 (10)	0.54892 (13)	0.0258 (6)
H38	0.5728	0.6949	0.5075	0.031*
C39	0.6650 (3)	0.69486 (10)	0.59258 (13)	0.0235 (6)
H39	0.7272	0.7095	0.5807	0.028*
C40	0.8220(3)	0 74817 (10)	0.69748(14)	0.0248(7)
C41	0.9313(3)	0.76130(11)	0.71415 (16)	0.0302(7)
H41	0.9865	0 7417	0.7312	0.036*
C42	0.9613(4)	0.80336(12)	0.70605(18)	0.0384(9)
H42	1.0362	0.8122	0.7186	0.046*
C43	0.8836(4)	0.8122 0.83164 (12)	0.68024 (19)	0.0479(11)
H43	0.9046	0.8600	0.6743	0.057*
C44	0.7730(4)	0.81873(11)	0.66248 (19)	0.037
H44	0.7189	0.8382	0.6439	0.058*
C45	0.7420(3)	0.0302 0.77722 (11)	0.67199 (16)	0.0354(8)
H45	0.6663	0.7687	0.6611	0.042*
Ασ?	0.224068 (16)	0.688496 (6)	0.323769 (9)	0.012 0.01375 (4)
Cl2	0.36640 (6)	0.65160(2)	0.40620 (3)	0.01616 (14)
N6	0.2600(2)	0 57945 (9)	0.30150(12)	0.0164 (5)
H6N	0.297(2)	0.6001 (8)	0.3217(14)	0.020*
N7	0.3116(2)	0.55402(9)	0.26416 (13)	0.0161 (5)
N8	0.0918(2)	0 56014 (9)	0 24624 (12)	0.0180(5)
H8N	0.129(3)	0 5404 (9)	0.2313(16)	0.022*
N9	0.5816(2)	0.51946 (9)	0.24039 (14)	0.0211 (6)
N10	0.4469 (2)	0.47841 (10)	0.14338 (13)	0.0218 (6)
P3	0.09628 (6)	0.74264 (3)	0.35680 (4)	0.01405(15)
P4	0.29648 (6)	0.69742(2)	0.22380 (3)	0.01487 (15)
S2	0.08349 (6)	0.62324(3)	0.32583 (4)	0.01625 (14)
C46	0.1472 (2)	0.58511(10)	0.28866 (14)	0.0144 (5)
C47	0.4171(2)	0.54722(10)	0 27876 (14)	0.0156 (6)
C48	0.4871(2)	0.56431(11)	0.33545 (15)	0.0188 (6)
H48A	0.4723	0.5944	0.3391	0.028*
H48B	0.5665	0 5600	0 3323	0.028*
H48C	0.4680	0.5496	0.3719	0.028*
C49	0.4690(2)	0.52074(10)	0.23486 (15)	0.0157 (6)
C50	0.6249(3)	0.49865(12)	0.19625(17)	0.0260(7)
H50	0 7040	0.4969	0 1986	0.031*
	0.7010	0.1202	0.1700	0.001

C51	0.5589 (3)	0.47956 (11)	0.14714 (16)	0.0233 (7)
H51	0.5938	0.4670	0.1154	0.028*
C52	0.4034 (3)	0.49866 (11)	0.18741 (15)	0.0185 (6)
Н52	0.3246	0.4982	0.1868	0.022*
C53	-0.0295(2)	0.56230(11)	0.22756 (15)	0.0228 (6)
H53A	-0.0479	0.5494	0.1863	0.027*
H53B	-0.0524	0.5922	0.2242	0.027*
C54	-0.0952(3)	0.54061 (12)	0.27121 (17)	0.0332 (8)
H54A	-0.1753	0.5434	0.2564	0.050*
H54B	-0.0784	0.5534	0.3121	0.050*
H54C	-0.0750	0.5108	0.2737	0.050*
C55	-0.0384(2)	0.73692 (9)	0.30721 (14)	0.0174 (6)
C56	-0.0344 (2)	0.72546 (10)	0.24668 (14)	0.0248 (6)
H56	0.0361	0.7206	0.2334	0.030*
C57	-0.1311(2)	0.72104 (11)	0.20526 (14)	0.0272 (7)
H57	-0.1264	0.7135	0.1638	0.033*
C58	-0.2340(2)	0.72749 (11)	0.22384 (14)	0.0266 (7)
H58	-0.3004	0.7248	0.1952	0.032*
C59	-0.2400(2)	0.73806 (10)	0.28518 (14)	0.0274 (7)
Н59	-0.3109	0.7421	0.2985	0.033*
C60	-0.1427 (2)	0.74268 (9)	0.32665 (13)	0.0217 (6)
H60	-0.1471	0.7498	0.3683	0.026*
C61	0.1279 (2)	0.79871 (10)	0.35066 (13)	0.0156 (6)
C62	0.0583 (2)	0.82645 (9)	0.31367 (13)	0.0181 (5)
H62	-0.0088	0.8165	0.2903	0.022*
C63	0.0874 (2)	0.86885 (9)	0.31104 (14)	0.0214 (6)
H63	0.0398	0.8878	0.2862	0.026*
C64	0.1865 (2)	0.88337 (10)	0.34501 (13)	0.0211 (6)
H64	0.2060	0.9123	0.3438	0.025*
C65	0.2558 (2)	0.85580 (9)	0.38011 (13)	0.0211 (6)
H65	0.3239	0.8657	0.4025	0.025*
C66	0.2272 (2)	0.81349 (9)	0.38338 (13)	0.0192 (5)
H66	0.2758	0.7947	0.4080	0.023*
C67	0.0663 (2)	0.73918 (10)	0.43656 (13)	0.0177 (6)
C68	0.0057 (2)	0.77009 (10)	0.46254 (14)	0.0230 (6)
H68	-0.0214	0.7938	0.4387	0.028*
C69	-0.0156 (2)	0.76655 (10)	0.52343 (14)	0.0254 (6)
H69	-0.0585	0.7873	0.5406	0.030*
C70	0.0269 (3)	0.73231 (11)	0.55850 (15)	0.0296 (7)
H70	0.0122	0.7294	0.5998	0.035*
C71	0.0904 (3)	0.70249 (12)	0.53351 (16)	0.0299 (8)
H71	0.1204	0.6794	0.5580	0.036*
C72	0.1110 (3)	0.70583 (10)	0.47272 (15)	0.0223 (6)
H72	0.1555	0.6853	0.4561	0.027*
C73	0.4135 (2)	0.66460 (10)	0.20850 (14)	0.0164 (6)
C74	0.5058 (2)	0.66061 (11)	0.25459 (14)	0.0210 (6)
H74	0.5047	0.6736	0.2935	0.025*
C75	0.5991 (3)	0.63762 (12)	0.24352 (16)	0.0244 (7)

H75	0.6627	0.6359	0.2743	0.029*
C76	0.5992 (3)	0.61724 (11)	0.18774 (16)	0.0229 (7)
H76	0.6630	0.6015	0.1804	0.028*
C77	0.5069 (3)	0.61954 (11)	0.14251 (15)	0.0218 (6)
H77	0.5067	0.6048	0.1048	0.026*
C78	0.4146(3)	0.64344(11)	0 15252 (15)	0.0180 (6)
H78	0.3518	0.6454	0.1212	0.022*
C79	0.3445(2)	0.75107(10)	0.1212 0.21133(13)	0.022
C80	0.3445(2) 0.4501(3)	0.75107(10) 0.76028(10)	0.21133(13) 0.10634(14)	0.0130(0)
	0.4501 (5)	0.70028 (10)	0.19034 (14)	0.0238 (0)
C 9 1	0.3009	0.7382 0.90220 (11)	0.1911 0.19902 (16)	0.029°
	0.4810 (5)	0.80230 (11)	0.18895 (10)	0.0290 (7)
H81	0.5544	0.8085	0.1/95	0.036*
C82	0.4080 (3)	0.83442 (11)	0.19526 (16)	0.0336 (8)
H82	0.4290	0.8627	0.1892	0.040*
C83	0.3028 (3)	0.82540 (10)	0.21054 (16)	0.0334 (8)
H83	0.2521	0.8476	0.2152	0.040*
C84	0.2712 (3)	0.78410 (9)	0.21902 (14)	0.0260 (6)
H84	0.1994	0.7782	0.2301	0.031*
C85	0.1917 (2)	0.68908 (10)	0.15654 (12)	0.0168 (5)
C86	0.1090 (2)	0.65863 (10)	0.15994 (13)	0.0208 (6)
H86	0.1085	0.6429	0.1968	0.025*
C87	0.0275 (3)	0.65125 (10)	0.10965 (14)	0.0248 (6)
H87	-0.0287	0.6307	0.1125	0.030*
C88	0.0278 (3)	0.67346 (11)	0.05595 (15)	0.0286 (8)
H88	-0.0277	0.6680	0.0217	0.034*
C89	0.1085 (3)	0.70359 (10)	0.05165 (13)	0.0277 (7)
H89	0.1081	0.7190	0.0145	0.033*
C90	0.1904 (2)	0.71145 (10)	0.10158 (14)	0.0242 (6)
H90	0.2460	0.7322	0.0983	0.029*
Ag3	0.473216 (16)	0.435776 (6)	0.576047 (9)	0.01369(5)
Cl3	0.61650 (6)	0 47296 (2)	0.65742(3)	0.01630(13)
N11	0.5127(2)	0.54412(9)	0.55109(12)	0.0157(5)
HIIN	0.5127(2) 0.547(2)	0.5309(10)	0.5837(11)	0.019*
N12	0.547(2)	0.56962 (0)	0.5037(11) 0.51280(12)	0.019
N12	0.3050(2)	0.56313(0)	0.31289(12) 0.40200(12)	0.0103(5)
HI3 HI2N	0.3403(2)	0.50313(9)	0.49290(12) 0.4736(15)	0.0175(3)
N14	0.381(3)	0.5821(9)	0.4730(13) 0.40304(14)	0.021°
IN 14	0.8304(2)	0.00499(9)	0.49304(14) 0.20420(12)	0.0224(6)
NI5	0.7056 (2)	0.04570(10)	0.39430(12)	0.0211(6)
P5	0.34543 (6)	0.38197 (3)	0.61001 (4)	0.01413(15)
P6	0.54480 (6)	0.42590 (3)	0.4/599 (3)	0.01462 (15)
83	0.33439 (6)	0.50170 (3)	0.57482 (4)	0.01697 (15)
C91	0.3997 (2)	0.53877 (10)	0.53675 (14)	0.0154 (6)
C92	0.6714 (2)	0.57670 (10)	0.52943 (14)	0.0159 (6)
C93	0.7392 (3)	0.56075 (12)	0.58698 (16)	0.0228 (7)
H93A	0.7154	0.5748	0.6227	0.034*
H93B	0.8186	0.5666	0.5857	0.034*
H93C	0.7280	0.5304	0.5903	0.034*
C94	0.7246 (2)	0.60256 (10)	0.48505 (14)	0.0158 (6)

C95	0.8817 (3)	0.62717 (12)	0.45030 (17)	0.0265 (8)
H95	0.9608	0.6296	0.4542	0.032*
C96	0.8178 (3)	0.64632 (12)	0.40133 (17)	0.0241 (7)
H96	0.8542	0.6604	0.3715	0.029*
C97	0.6597 (3)	0.62427 (11)	0.43676 (15)	0.0180 (6)
H97	0.5805	0.6237	0.4343	0.022*
C98	0.2258 (2)	0.56039 (11)	0.47262 (15)	0.0213 (6)
H98A	0.2033	0.5304	0.4695	0.026*
H98B	0.2084	0.5731	0.4311	0.026*
C99	0.1591 (3)	0.58266 (13)	0.51643 (18)	0.0378 (9)
H99A	0.0790	0.5798	0.5016	0.057*
H99B	0.1794	0.6125	0.5186	0.057*
H99C	0.1756	0.5701	0.5575	0.057*
C100	0.3150 (2)	0.38481 (9)	0.68954 (13)	0.0173 (6)
C101	0.2538 (2)	0.35355 (9)	0.71551 (13)	0.0219 (6)
H101	0.2265	0.3300	0.6913	0.026*
C102	0.2326 (2)	0.35663 (11)	0.77614 (14)	0.0258 (6)
H102	0.1904	0.3355	0.7931	0.031*
C103	0.2737 (3)	0.39084 (10)	0.81157 (14)	0.0282 (7)
H103	0.2580	0.3937	0.8526	0.034*
C104	0.3379 (3)	0.42092 (12)	0.78689 (16)	0.0321 (8)
H104	0.3684	0.4436	0.8119	0.039*
C105	0.3584 (3)	0.41839 (10)	0.72604 (15)	0.0248 (7)
H105	0.4017	0.4394	0.7096	0.030*
C106	0.3779 (2)	0.32625 (10)	0.60382 (14)	0.0154 (6)
C107	0.4774 (2)	0.31149 (9)	0.63677 (13)	0.0202 (5)
H107	0.5256	0.3304	0.6614	0.024*
C108	0.5068 (2)	0.26930 (10)	0.63396 (13)	0.0228 (6)
H108	0.5756	0.2597	0.6560	0.027*
C109	0.4364 (3)	0.24128 (10)	0.59929 (14)	0.0214 (6)
H109	0.4555	0.2123	0.5985	0.026*
C110	0.3367 (2)	0.25581 (9)	0.56538 (13)	0.0196 (6)
H110	0.2889	0.2368	0.5408	0.023*
C111	0.3075 (2)	0.29808 (9)	0.56751 (13)	0.0185 (5)
H111	0.2400	0.3078	0.5444	0.022*
C112	0.2114 (2)	0.38766 (9)	0.56068 (13)	0.0165 (5)
C113	0.1067(2)	0.38141 (9)	0.57933 (13)	0.0207(5)
H113	0.1017	0.3740	0.6208	0.025*
C114	0.0087(2)	0 38603 (10)	0 53726 (14)	0.0250 (6)
H114	-0.0623	0.3818	0.5503	0.030*
C115	0.0154 (2)	0.39675 (11)	0.47695 (15)	0.0268 (7)
H115		0 3995	0 4484	0.032*
C116	0.1180 (3)	0.40352 (11)	0.45798 (14)	0.0287(7)
H116	0 1224	0.4112	0.4165	0.034*
C117	0.2154 (2)	0.39908 (10)	0.49974 (13)	0.0235 (6)
H117	0.2860	0.4039	0.4864	0.028*
C118	0.6609 (2)	0.45917 (10)	0.45967 (14)	0.0170 (6)
C119	0.7531 (2)	0.46313 (11)	0.50541 (15)	0.0207 (6)

H119	0.7527	0.4499	0.5442	0.025*
C120	0.8460 (3)	0.48661 (12)	0.49410 (16)	0.0228 (7)
H120	0.9090	0.4891	0.5252	0.027*
C121	0.8467 (3)	0.50637 (11)	0.43751 (16)	0.0221 (7)
H121	0.9106	0.5217	0.4295	0.027*
C122	0.7533 (3)	0.50349 (11)	0.39270 (15)	0.0219 (6)
H122	0.7526	0.5176	0.3545	0.026*
C123	0.6604 (3)	0.47990 (10)	0.40375 (15)	0.0182 (6)
H123	0.5968	0.4780	0.3730	0.022*
C124	0.4400 (2)	0.43286 (10)	0.40783 (12)	0.0172 (5)
C125	0.4387 (2)	0.40877 (10)	0.35524 (13)	0.0227 (6)
H125	0.4932	0.3874	0.3538	0.027*
C126	0.35831 (19)	0.41567 (8)	0.30446 (10)	0.0281 (7)
H126	0.3575	0.3987	0.2687	0.034*
C127	0.27943 (19)	0.44705 (8)	0.30562 (10)	0.0266 (7)
H127	0.2253	0.4520	0.2705	0.032*
C128	0.2796 (3)	0.47113 (11)	0.35795 (15)	0.0271 (7)
H128	0.2253	0.4926	0.3590	0.032*
C129	0.3596 (3)	0.46393 (10)	0.40927 (14)	0.0223 (6)
H129	0.3591	0.4804	0.4454	0.027*
C130	0.5952 (2)	0.37254 (10)	0.46526 (13)	0.0186 (6)
C131	0.7004 (3)	0.36319 (11)	0.45056 (15)	0.0241 (7)
H131	0.7507	0.3854	0.4452	0.029*
C132	0.7338 (3)	0.32179 (12)	0.44343 (17)	0.0341 (8)
H132	0.8064	0.3160	0.4334	0.041*
C133	0.6611 (3)	0.28918 (11)	0.45095 (16)	0.0344 (8)
H133	0.6835	0.2609	0.4456	0.041*
C134	0.5567 (3)	0.29753 (10)	0.46617 (15)	0.0321 (7)
H134	0.5071	0.2751	0.4717	0.039*
C135	0.5238 (3)	0.33892 (9)	0.47343 (13)	0.0255 (6)
H135	0.4515	0.3445	0.4841	0.031*
Ag4	-0.038125 (18)	0.439474 (6)	0.073147 (9)	0.01726 (5)
Cl4	0.10740 (6)	0.47297 (2)	0.15720 (3)	0.01678 (14)
N16	0.0143 (2)	0.54561 (9)	0.05256 (13)	0.0156 (5)
H16N	0.047 (3)	0.5308 (9)	0.0825 (12)	0.019*
N17	0.0680 (2)	0.57030 (9)	0.01351 (12)	0.0154 (5)
N18	-0.1502 (2)	0.56879 (10)	-0.00276 (12)	0.0192 (5)
H18N	-0.112 (3)	0.5859 (10)	-0.0222 (16)	0.023*
N19	0.3403 (2)	0.59985 (10)	-0.01125 (14)	0.0248 (6)
N20	0.2088 (2)	0.64353 (10)	-0.10746 (13)	0.0236 (6)
P7	-0.15849 (6)	0.38179 (3)	0.10226 (4)	0.01484 (15)
P8	0.01543 (7)	0.43194 (3)	-0.03167 (3)	0.02084 (16)
S4	-0.16819 (6)	0.50928 (3)	0.08195 (4)	0.01785 (15)
C136	-0.0996 (2)	0.54359 (10)	0.04060 (14)	0.0160 (6)
C137	0.1748 (2)	0.57632 (10)	0.02937 (14)	0.0157 (6)
C138	0.2424 (3)	0.56014 (12)	0.08687 (15)	0.0197 (7)
H13A	0.2202	0.5748	0.1226	0.030*
H13B	0.3220	0.5652	0.0851	0.030*

H13C	0.2293	0.5299	0.0906	0.030*
C139	0.2285 (2)	0.60097 (10)	-0.01585 (14)	0.0162 (6)
C140	0.3846 (3)	0.61961 (14)	-0.05631 (18)	0.0346 (9)
H140	0.4636	0.6195	-0.0552	0.041*
C141	0.3205 (3)	0.63988 (13)	-0.10371 (18)	0.0310 (8)
H141	0.3566	0.6519	-0.1354	0.037*
C142	0.1638 (3)	0.62441 (11)	-0.06249 (14)	0.0182 (6)
H142	0.0854	0.6267	-0.0619	0.022*
C143	-0.2710 (3)	0.57430 (14)	-0.01568 (17)	0.0388 (10)
H14A	-0.3090	0.5475	-0.0087	0.047*
H14B	-0.2924	0.5825	-0.0593	0.047*
C144	-0.3074 (4)	0.60825 (19)	0.0264 (2)	0.0759 (19)
H14C	-0.3887	0.6119	0.0179	0.114*
H14D	-0.2704	0.6348	0.0189	0.114*
H14E	-0.2863	0.5999	0.0695	0.114*
C145	-0.2137 (2)	0.39298 (10)	0.17381 (13)	0.0212 (6)
C146	-0.3248 (3)	0.38613 (9)	0.18182 (14)	0.0252 (6)
H146	-0.3767	0.3760	0.1485	0.030*
C147	-0.3604 (3)	0.39411 (10)	0.23859 (16)	0.0330(7)
H147	-0.4362	0.3895	0.2439	0.040*
C148	-0.2845 (3)	0.40885 (11)	0.28712 (16)	0.0382 (8)
H148	-0.3082	0.4137	0.3261	0.046*
C149	-0.1737 (4)	0.41671 (13)	0.27939 (16)	0.0335 (8)
H149	-0.1223	0.4270	0.3129	0.040*
C150	-0.1386 (3)	0.40949 (11)	0.22258 (16)	0.0278 (7)
H150	-0.0638	0.4157	0.2167	0.033*
C151	-0.2832 (2)	0.37740 (10)	0.04449 (14)	0.0195 (6)
C152	-0.3356 (2)	0.41434 (9)	0.02134 (13)	0.0217 (6)
H152	-0.3062	0.4407	0.0360	0.026*
C153	-0.4306 (3)	0.41300 (11)	-0.02306 (15)	0.0274 (7)
H153	-0.4657	0.4384	-0.0380	0.033*
C154	-0.4739 (3)	0.37496 (11)	-0.04535 (15)	0.0297 (7)
H154	-0.5388	0.3741	-0.0754	0.036*
C155	-0.4219 (3)	0.33781 (11)	-0.02347 (15)	0.0325 (7)
H155	-0.4510	0.3115	-0.0388	0.039*
C156	-0.3271 (3)	0.33913 (10)	0.02085 (14)	0.0265 (6)
H156	-0.2917	0.3136	0.0352	0.032*
C157	-0.1113 (2)	0.32737 (10)	0.11259 (14)	0.0170 (6)
C158	-0.1672 (2)	0.29837 (9)	0.14576 (13)	0.0229 (6)
H158	-0.2268	0.3075	0.1666	0.027*
C159	-0.1352 (2)	0.25615 (9)	0.14803 (14)	0.0240 (6)
H159	-0.1741	0.2365	0.1698	0.029*
C160	-0.0466 (3)	0.24285 (10)	0.11850 (14)	0.0236 (6)
H160	-0.0258	0.2140	0.1194	0.028*
C161	0.0111 (2)	0.27170 (10)	0.08788 (14)	0.0258 (6)
H161	0.0729	0.2627	0.0687	0.031*
C162	-0.0205 (2)	0.31398 (9)	0.08476 (14)	0.0224 (6)
H162	0.0199	0.3336	0.0637	0.027*

C163	0.1388 (3)	0.46079 (10)	-0.04771 (14)	0.0187 (6)
C164	0.2273 (3)	0.46542 (10)	-0.00002 (14)	0.0219 (6)
H164	0.2201	0.4546	0.0397	0.026*
C165	0.3260 (3)	0.48554 (12)	-0.00943 (16)	0.0249 (7)
H165	0.3863	0.4878	0.0232	0.030*
C166	0.3354 (3)	0.50225 (12)	-0.06692 (16)	0.0257 (7)
H166	0.4022	0.5162	-0.0736	0.031*
C167	0.2474 (3)	0.49856 (11)	-0.11457 (15)	0.0242 (7)
H167	0.2540	0.5102	-0.1538	0.029*
C168	0.1494 (3)	0.47781 (11)	-0.10522 (15)	0.0209 (6)
H168	0.0896	0.4753	-0.1381	0.025*
C169	0.0499 (3)	0.37773 (10)	-0.04887 (14)	0.0304 (7)
C170	-0.0360 (4)	0.34958 (11)	-0.07160 (16)	0.0453 (10)
H170	-0.1112	0.3591	-0.0804	0.054*
C171	-0.0100 (5)	0.30719 (12)	-0.08127 (18)	0.0650 (15)
H171	-0.0672	0.2883	-0.0981	0.078*
C172	0.0998 (5)	0.29289 (13)	-0.06615 (19)	0.0625 (15)
H172	0.1168	0.2641	-0.0714	0.075*
C173	0.1832 (4)	0.32030 (12)	-0.04381 (19)	0.0507 (11)
H173	0.2580	0.3105	-0.0343	0.061*
C174	0.1590 (4)	0.36288 (11)	-0.03481 (16)	0.0367 (9)
H174	0.2174	0.3816	-0.0191	0.044*
C175	-0.0980 (3)	0.44532 (11)	-0.09379 (13)	0.0245 (7)
C176	-0.1838 (3)	0.47031 (11)	-0.07890 (14)	0.0286 (7)
H176	-0.1809	0.4807	-0.0381	0.034*
C177	-0.2757 (3)	0.48059 (11)	-0.12342 (15)	0.0324 (7)
H177	-0.3336	0.4982	-0.1129	0.039*
C178	-0.2808 (3)	0.46484 (11)	-0.18230 (14)	0.0312 (7)
H178	-0.3435	0.4709	-0.2123	0.037*
C179	-0.1953 (3)	0.44043 (11)	-0.19746 (14)	0.0295 (7)
H179	-0.1987	0.4302	-0.2384	0.035*
C180	-0.1033 (3)	0.43022 (11)	-0.15419 (14)	0.0302 (7)
H180	-0.0448	0.4132	-0.1655	0.036*

Atomic displacement parameters $(Å^2)$

	U^{11}	U^{22}	U^{33}	U^{12}	U^{13}	U^{23}
Agl	0.01779 (9)	0.01646 (10)	0.01339 (9)	0.00323 (8)	0.00165 (7)	0.00037 (8)
Cl1	0.0175 (3)	0.0188 (4)	0.0146 (3)	0.0009 (3)	-0.0022 (3)	-0.0005 (3)
N1	0.0170 (11)	0.0185 (14)	0.0177 (13)	-0.0021 (10)	0.0020 (10)	-0.0071 (11)
N2	0.0189 (12)	0.0146 (13)	0.0182 (13)	0.0016 (10)	0.0030 (11)	-0.0031 (11)
N3	0.0180 (11)	0.0188 (13)	0.0189 (13)	-0.0015 (10)	0.0033 (10)	-0.0044 (11)
N4	0.0160 (12)	0.0261 (15)	0.0267 (14)	0.0014 (10)	0.0008 (11)	-0.0082 (12)
N5	0.0227 (12)	0.0221 (14)	0.0207 (13)	-0.0002 (11)	0.0043 (11)	-0.0037 (11)
P1	0.0162 (3)	0.0129 (4)	0.0157 (3)	0.0015 (3)	0.0027 (3)	-0.0005 (3)
P2	0.0239 (3)	0.0173 (4)	0.0134 (3)	0.0028 (3)	0.0047 (3)	0.0016 (3)
S 1	0.0149 (3)	0.0186 (3)	0.0161 (3)	-0.0003 (3)	0.0040 (3)	-0.0017 (3)
C1	0.0181 (13)	0.0131 (13)	0.0129 (13)	-0.0019 (10)	0.0022 (11)	-0.0010 (11)

C2	0.0158 (12)	0.0159 (15)	0.0157 (14)	-0.0009 (11)	-0.0009 (11)	0.0007 (11)
C3	0.0216 (15)	0.0210 (17)	0.0220 (16)	0.0021 (13)	-0.0007 (13)	-0.0021 (14)
C4	0.0151 (12)	0.0137 (14)	0.0199 (14)	0.0012 (10)	0.0027 (11)	-0.0017 (11)
C5	0.0170 (14)	0.0338 (18)	0.0315 (18)	0.0025 (13)	0.0020 (13)	-0.0143 (15)
C6	0.0223 (14)	0.0251 (16)	0.0238 (16)	0.0025 (12)	0.0039 (12)	-0.0090 (13)
C7	0.0178 (13)	0.0179 (15)	0.0188 (15)	-0.0019(12)	0.0021 (12)	-0.0017(12)
C8	0.0163 (13)	0.0357 (17)	0.0212 (14)	-0.0064(12)	0.0018 (11)	-0.0073(13)
C9	0.0327 (17)	0.049 (2)	0.0318 (18)	-0.0251 (16)	0.0083 (14)	-0.0037(16)
C10	0.0194(13)	0.0148 (14)	0.0144 (13)	0.0006 (10)	0.0008 (10)	-0.0001(11)
C11	0.0211 (13)	0.0227 (14)	0.0177 (13)	-0.0011(11)	0.0024 (11)	0.0014 (11)
C12	0.0239(14)	0.0249(15)	0.0199 (14)	-0.0057(11)	0.0011 (11)	0.0011 (11)
C13	0.0319 (16)	0.0192(15)	0.0216 (15)	-0.0076(12)	0.0014(12)	-0.0014(12)
C14	0.0322(15)	0.0192(12) 0.0191(14)	0.0263(15)	0.0031(12)	0.0043(12)	-0.0040(12)
C15	0.0322(13) 0.0234(13)	0.0211(14)	0.0200(12) 0.0240(14)	0.0001(12)	0.0013(12) 0.0062(11)	-0.0002(12)
C16	0.0238(13)	0.0211(11) 0.0117(13)	0.0210(11) 0.0202(14)	0.00000(11) 0.0026(10)	0.0002(11) 0.0078(11)	0.0002(12)
C17	0.0290(15)	0.0228(14)	0.0202(11) 0.0252(15)	0.0020(10) 0.0012(12)	0.0070(11) 0.0061(12)	0.0016(11) 0.0026(12)
C18	0.0290(19) 0.0385(18)	0.0220(11) 0.0292(17)	0.0232(13) 0.0337(18)	0.0012(12) 0.0024(14)	0.0001(12) 0.0202(15)	0.0020(12) 0.0028(14)
C10	0.0505(10)	0.0292(17) 0.0206(15)	0.0308(18)	0.0024(14)	0.0202(13) 0.0247(17)	0.0020(14) 0.0053(13)
C20	0.002(2) 0.052(2)	0.0200(13) 0.045(2)	0.0203(17)	0.0090(18)	0.0247(17)	0.0055(15)
C20	0.032(2)	0.045(2)	0.0205(17) 0.0185(15)	0.0109(10)	0.0030(10) 0.0041(13)	0.0000(10)
C21	0.0311(10) 0.0177(12)	0.043(2) 0.0222(14)	0.0163(13)	0.0074(13) 0.0036(11)	0.0041(10) 0.0025(10)	0.0009(13) 0.0004(11)
C23	0.0177(12) 0.0335(16)	0.0222(11) 0.0195(14)	0.0278(16)	0.0000(11) 0.0002(12)	-0.0097(13)	-0.0017(12)
C24	0.0394(18)	0.0155(11)	0.0270(10) 0.0344(18)	0.0002(12) 0.0061(14)	-0.0158(14)	0.0017(12)
C25	0.0354(10) 0.0253(15)	0.0235(10) 0.0385(19)	0.0270(17)	0.0001(14)	-0.0068(12)	-0.0025(14)
C26	0.0233(15) 0.0284(15)	0.0305(15) 0.0246(16)	0.0270(17) 0.0280(17)	-0.0037(13)	-0.0012(13)	-0.0007(13)
C27	0.0201(13)	0.0210(10) 0.0223(14)	0.0200(17) 0.0273(16)	0.0000(11)	0.0012(13)	0.00007(13)
C28	0.0210(13) 0.0223(13)	0.0223(11) 0.0150(14)	0.0279(10) 0.0189(14)	-0.0012(11)	0.0017(12) 0.0067(11)	0.0008(12)
C29	0.0223(13) 0.0203(14)	0.0130(11) 0.0238(16)	0.0103(14)	-0.0012(11)	0.0007(11) 0.0045(11)	0.0000(11)
C30	0.0203(11) 0.0248(14)	0.0258(16)	0.0175(11) 0.0205(14)	0.0022(12)	0.0015(11) 0.0085(12)	-0.0039(12)
C31	0.0210(14)	0.0283(17)	0.0251 (16)	-0.0001(13)	0.00000(12)	-0.0014(13)
C32	0.0227(15)	0.0203(17) 0.0327(19)	0.0222(16)	0.0005(13)	0.0071(13)	0 0006 (14)
C33	0.0220(15)	0.0209(15)	0.0195(14)	-0.0033(12)	0.0036(12)	-0.0023(12)
C34	0.0284(14)	0.0210(16)	0.0136(12)	0.00000(12)	0.0030(12)	0.0023(12)
C35	0.0207(11) 0.0302(15)	0.0246(17)	0.0130(12) 0.0139(13)	-0.0010(13)	-0.0010(12)	0.0055(11)
C36	0.0202(15) 0.0292(16)	0.0310(17)	0.0228 (16)	-0.0017(15)	-0.0006(13)	0.0002(12) 0.0036(15)
C37	0.0366(17)	0.0274(16)	0.0182(14)	0.0045(13)	-0.0022(13)	-0.0033(12)
C38	0.0300(17) 0.0416(17)	0.0234(15)	0.0102(11) 0.0117(12)	0.0016(13)	0.0011(11)	0.0019(12)
C39	0.0340(15)	0.0231(15)	0.0135(12)	0.00000(11)	0.0040(11)	0.0019(12)
C40	0.0310(13) 0.0419(17)	0.0231(13) 0.0172(14)	0.0155(12) 0.0164(14)	0.0002(12)	0.0010(11) 0.0080(13)	0.0000(11) 0.0026(11)
C41	0.0439(19)	0.0172(11) 0.0209(16)	0.0101(11) 0.0247(17)	-0.0029(12)	0.0000(13) 0.0010(14)	0.0020(11) 0.0021(13)
C42	0.058(2)	0.0249(18)	0.0217(17) 0.0312(19)	-0.0120(16)	0.0010(11) 0.0024(16)	0.0021(15)
C43	0.033(2) 0.087(3)	0.0219(10) 0.0208(17)	0.036(2)	-0.0072(19)	0.0021(10)	0.0002(10)
C44	0.081(3)	0.0213(17)	0.040(2)	0.0112(18)	-0.003(2)	0.0015 (16)
C45	0.0469 (19)	0.0273 (16)	0.0301 (17)	0.0072(15)	-0.0009(15)	0.0011 (14)
Ag2	0.01437 (9)	0.01432 (10)	0.01253 (10)	0.00230 (8)	0.00178 (7)	0.00075 (8)
Cl2	0.0160 (3)	0.0171 (3)	0.0142 (3)	0.0014 (3)	-0.0021(2)	0.0004 (3)
N6	0.0168 (11)	0.0171 (13)	0.0149 (12)	0.0008 (10)	0.0014 (10)	-0.0046 (10)
N7	0.0173 (11)	0.0145 (13)	0.0171 (13)	0.0021 (10)	0.0040 (10)	-0.0014 (10)
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N8	0.0177(11)	0.0178(12)	0.0185 (12)	-0.0010(10)	0.0022 (10)	-0.0054(10)
N9	0.0174(11)	0.0211(13)	0.0245(14)	0.0025(10)	0.0024(10)	-0.0042(11)
N10	0.0246(13)	0.0214(14)	0.0199(13)	0.0020(10)	0.0053(11)	-0.0022(11)
P3	0.0139(3)	0.0138(4)	0.0148(3)	0.0011 (3)	0.0029 (3)	0.0005(3)
P4	0.0161(3)	0.0169 (4)	0.0121(3)	0.0012(3)	0.0035(2)	0.0012(3)
S2	0.0101(3) 0.0127(3)	0.0180(3)	0.0121(3) 0.0184(3)	-0.0007(2)	0.0035(2)	-0.0040(3)
C46	0.0127(3) 0.0157(12)	0.0130(3)	0.0101(3) 0.0145(13)	-0.0018(10)	0.0023(11)	0.0000(11)
C47	0.0175(13)	0.0165(15)	0.0122(13)	-0.0009(11)	-0.0003(11)	-0.0006(11)
C48	0.0179(13) 0.0168(13)	0.0103 (15)	0.0122(15) 0.0198(15)	0.0009(11) 0.0027(12)	-0.0002(12)	-0.0023(13)
C49	0.0159(12)	0.0125(13)	0.0198(13)	0.0027(12) 0.0007(10)	0.0002(12) 0.0030(11)	0.0025(15)
C50	0.0201(12)	0.0123(13) 0.0274(17)	0.0313(18)	0.0007(10) 0.0031(13)	0.0058(13)	-0.0062(14)
C51	0.0201(11) 0.0247(15)	0.0238(16)	0.0211(15)	0.0031(13) 0.0041(12)	0.0055(12)	-0.0066(13)
C52	0.0217(13)	0.0230(10) 0.0184(15)	0.0221(13) 0.0179(14)	0.0011(12) 0.0038(12)	0.0033(12) 0.0017(11)	-0.0011(12)
C53	0.0191(13) 0.0164(12)	0.0296 (16)	0.0175(14)	-0.0039(11)	-0.0003(11)	-0.0046(12)
C54	0.0101(12) 0.0277(16)	0.0290(10)	0.0217(11) 0.0324(18)	-0.0145(14)	0.0003 (11)	0.0010(12)
C55	0.0277(10) 0.0143(12)	0.0169(14)	0.0327(10)	0.003(10)	0.0014(10)	-0.0001(13)
C56	0.0143(12) 0.0217(14)	0.0109(14) 0.0298(16)	0.0207(14) 0.0229(15)	0.0003(10) 0.0001(12)	0.0014(10) 0.0038(11)	-0.0006(11)
C57	0.0219(14)	0.0293(10) 0.0407(19)	0.0229(13) 0.0186(14)	0.0001(12) 0.0027(13)	0.0030(11) 0.0011(11)	-0.007(13)
C58	0.0219(14) 0.0183(13)	0.0361(18)	0.0100(14) 0.0240(15)	-0.0027(13)	-0.0018(12)	-0.0027(13)
C59	0.0165(13)	0.0364(18)	0.0240(15) 0.0302(16)	0.0009(13)	0.0010(12)	-0.0023(13)
C60	0.0105(13) 0.0198(13)	0.0268(15)	0.0302(10) 0.0191(13)	0.0010(12) 0.0020(11)	0.0000(12) 0.0043(10)	-0.0015(11)
C61	0.0190(12)	0.0200(12) 0.0173(14)	0.0143(12)	0.0020(11) 0.0003(10)	0.0073(10)	-0.0029(11)
C62	0.0166(12)	0.0177(13)	0.0203(13)	0.0003(10)	0.0075(10)	-0.0013(10)
C63	0.0217(14)	0.0197(14)	0.0230(13)	0.0003(10) 0.0044(11)	0.0036(11)	0.0019(10)
C64	0.0217(11) 0.0277(14)	0.0197(11) 0.0182(14)	0.0230(14)	-0.0042(12)	0.0030(11) 0.0084(12)	-0.0024(12)
C65	0.0279(13)	0.0102(11) 0.0219(14)	0.0190(11) 0.0189(13)	-0.0064(11)	0.0001(12) 0.0043(11)	-0.0027(11)
C66	0.0229(13)	0.0219(11) 0.0206(13)	0.0109(13) 0.0170(13)	0.0006 (10)	0.0013(11) 0.0023(10)	0.0027(11)
C67	0.0200(13) 0.0186(13)	0.0200(13) 0.0178(14)	0.0178(13)	0.0008(11)	0.0023(10) 0.0068(11)	0.0000(11)
C68	0.0224(13)	0.0250(15)	0.0217(14)	0.0023(11)	0.0031(11)	-0.0025(12)
C69	0.0245(14)	0.0316 (16)	0.0209(14)	0.0002(12)	0.0066 (11)	-0.0077(12)
C70	0.0213(11) 0.0361(17)	0.0336(18)	0.0209(11) 0.0204(15)	-0.0071(14)	0.0000(11) 0.0092(13)	0.000(13)
C71	0.047(2)	0.0216 (16)	0.0229 (16)	-0.0007(14)	0.0095(15)	0.0050(13)
C72	0.0305(15)	0.0185(15)	0.0183(14)	0.0003(12)	0.0048 (12)	0.0028(11)
C73	0.0185 (12)	0.0151 (13)	0.0159 (13)	0.0002 (10)	0.0042 (10)	0.0016 (11)
C74	0.0226(14)	0.0252(16)	0.0145 (13)	0.0031(12)	0.0006(11)	-0.0004(12)
C75	0.0210 (14)	0.0313 (18)	0.0203 (15)	0.0046 (13)	0.0010 (12)	0.0022 (13)
C76	0.0206 (14)	0.0257 (16)	0.0236(15)	0.0046 (12)	0.0068 (12)	0.0017(13)
C77	0.0241(14)	0.0232(15)	0.0194 (14)	0.0026(12)	0.0076(12)	-0.0033(12)
C78	0.0217(11) 0.0167(12)	0.0232(15)	0.0150(13)	0.0020(12) 0.0004(11)	0.0079(12)	0.00000(12)
C79	0.0217(13)	0.0179 (13)	0.0140(12)	-0.0027(11)	0.0007(10)	0.0010 (10)
C80	0.0250(14)	0.0251 (16)	0.0201(14)	-0.0030(12)	-0.0009(11)	0.0009(12)
C81	0.0266 (15)	0.0302(17)	0.0316 (17)	-0.0097(13)	0.0027 (13)	0.0051 (14)
C82	0.055 (2)	0.0216 (16)	0.0242 (16)	-0.0104(15)	0.0056 (15)	0.0041 (13)
C83	0.053(2)	0.0186 (15)	0.0311(17)	0.0067 (14)	0.0128 (16)	0.0019 (13)
C84	0.0323 (16)	0.0197 (14)	0.0281 (16)	0.0016 (12)	0.0119 (13)	0.0034 (12)
C85	0.0199 (12)	0.0189 (13)	0.0113 (11)	0.0065 (11)	0.0013 (9)	-0.0004(10)
C86	0.0267 (14)	0.0200 (14)	0.0144 (13)	0.0018 (12)	-0.0014 (11)	-0.0011 (11)
C87	0.0276 (15)	0.0234 (15)	0.0222 (15)	-0.0021(12)	-0.0013(12)	-0.0046(12)
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C88	0.0253 (15)	0.0383 (19)	0.0202 (15)	0.0105 (14)	-0.0038 (12)	-0.0117 (13)
C89	0.0340 (16)	0.0348 (17)	0.0140 (13)	0.0083 (13)	0.0017 (12)	0.0052 (12)
C90	0.0229 (14)	0.0332 (16)	0.0175 (13)	0.0056 (12)	0.0060 (11)	0.0053 (12)
Ag3	0.01369 (9)	0.01420 (12)	0.01306 (9)	-0.00204 (8)	0.00143 (7)	-0.00095 (8)
C13	0.0163 (3)	0.0168 (3)	0.0146 (3)	-0.0025 (3)	-0.0019 (3)	0.0001 (3)
N11	0.0162 (11)	0.0147 (12)	0.0159 (12)	-0.0026 (9)	0.0015 (10)	0.0030 (10)
N12	0.0153 (11)	0.0168 (13)	0.0170 (12)	-0.0005 (10)	0.0027 (10)	0.0012 (11)
N13	0.0148 (11)	0.0179 (13)	0.0195 (13)	0.0023 (9)	0.0018 (10)	0.0052 (10)
N14	0.0162 (11)	0.0238 (15)	0.0269 (15)	-0.0017 (10)	0.0018 (11)	0.0066 (12)
N15	0.0203 (12)	0.0250 (15)	0.0177 (13)	0.0005 (10)	0.0009 (10)	0.0048 (11)
P5	0.0130 (3)	0.0142 (4)	0.0153 (3)	-0.0018(3)	0.0025 (3)	-0.0001(3)
P6	0.0151 (3)	0.0165 (4)	0.0125 (3)	-0.0019(3)	0.0028 (2)	-0.0012(3)
S 3	0.0148 (3)	0.0170 (4)	0.0196 (4)	0.0010 (3)	0.0042 (3)	0.0038 (3)
C91	0.0148(12)	0.0154 (14)	0.0167 (14)	-0.0001(11)	0.0046(11)	-0.0021(11)
C92	0.0189(13)	0.0110 (13)	0.0180 (14)	-0.0029(11)	0.0032(11)	0.0003(11)
C93	0.0203(14)	0.0250(18)	0.0215(16)	-0.0023(13)	-0.0023(13)	0.00002(11) 0.0100(14)
C94	0.0200(11) 0.0180(13)	0.0250(10) 0.0150(14)	0.0212(13)	-0.0015(11)	0.0023(13)	-0.0022(11)
C95	0.0100(13) 0.0177(14)	0.0130(11) 0.0287(18)	0.0112(13) 0.0324(18)	-0.0019(11)	0.0012(11) 0.0006(13)	0.0022(11)
C96	0.0195(14)	0.0207(18)	0.0321(10) 0.0270(17)	-0.0020(13)	0.0000(13) 0.0074(13)	0.0110(13) 0.0054(14)
C97	0.0175(13)	0.0270(10) 0.0194(15)	0.0270(17) 0.0163(14)	-0.0007(12)	-0.0004(11)	0.0094(14)
C98	0.0175(13)	0.0194(15) 0.0244(16)	0.0103(14) 0.0204(14)	0.0007(11) 0.0041(11)	0.0004(11)	0.0009(12) 0.0048(12)
C90	0.0100(13) 0.0298(17)	0.0244(10)	0.0204(14)	0.0041(11) 0.0198(16)	0.0010(11) 0.0078(15)	0.0048(12)
C100	0.0298(17)	0.044(2)	0.040(2)	0.0130(10)	0.0078(13)	-0.0009(18)
C100	0.0190(13)	0.0171(13)	0.0137(13)	-0.0030(11)	0.0045(11)	-0.0002(11)
C101	0.0215(13)	0.0232(14) 0.0344(17)	0.0210(14) 0.0227(15)	-0.0022(11)	0.0040(11) 0.0074(11)	-0.0001(11)
C102	0.0213(14)	0.0344(17)	0.0227(13)	0.0004(12)	0.0074(11)	0.0077(13)
C105	0.05/4(17)	0.0312(17)	0.0184(14)	0.0000(14)	0.0120(15)	0.0013(12)
C104	0.051(2)	0.024/(17)	0.0219 (16)	-0.0029(13)	0.0095(15)	-0.0040(13)
C105	0.0341(16)	0.0200 (15)	0.0218(15)	-0.0058(12)	0.0092(13)	-0.0006(12)
C106	0.0164 (12)	0.0141 (13)	0.0168 (13)	-0.0002(10)	0.0060 (10)	-0.0014 (11)
C107	0.0208 (13)	0.0232 (14)	0.0162 (13)	-0.0006 (11)	0.0011 (10)	-0.0013 (11)
C108	0.0235 (13)	0.0264 (15)	0.0179 (13)	0.0039 (12)	0.0012 (11)	0.0026 (12)
C109	0.0306 (15)	0.0140 (13)	0.0218 (14)	0.0032 (12)	0.0115 (12)	0.0012 (11)
C110	0.0241 (13)	0.0162 (13)	0.0197 (13)	-0.0005 (11)	0.0077 (11)	-0.0023 (11)
C111	0.0162 (12)	0.0198 (13)	0.0195 (13)	-0.0011 (10)	0.0030 (10)	-0.0023 (11)
C112	0.0145 (11)	0.0156 (13)	0.0190 (13)	-0.0010 (10)	0.0004 (10)	0.0000 (10)
C113	0.0179 (12)	0.0217 (13)	0.0229 (14)	-0.0012 (11)	0.0045 (10)	0.0050 (11)
C114	0.0154 (13)	0.0308 (16)	0.0292 (15)	0.0008 (11)	0.0048 (11)	0.0032 (13)
C115	0.0198 (14)	0.0330 (17)	0.0260 (16)	0.0016 (12)	-0.0028 (12)	0.0003 (13)
C116	0.0261 (15)	0.0401 (19)	0.0186 (14)	-0.0029 (14)	-0.0011 (12)	0.0042 (13)
C117	0.0163 (13)	0.0332 (16)	0.0210 (14)	-0.0032 (12)	0.0024 (11)	0.0026 (12)
C118	0.0164 (12)	0.0185 (14)	0.0165 (13)	-0.0009 (11)	0.0038 (10)	-0.0046 (11)
C119	0.0233 (14)	0.0226 (16)	0.0164 (14)	-0.0038 (12)	0.0037 (11)	-0.0018 (12)
C120	0.0211 (14)	0.0270 (17)	0.0193 (15)	-0.0048 (12)	0.0000 (12)	-0.0007 (13)
C121	0.0180 (13)	0.0238 (16)	0.0254 (16)	-0.0052 (12)	0.0060 (12)	-0.0024 (13)
C122	0.0219 (14)	0.0235 (16)	0.0212 (15)	-0.0033 (12)	0.0056 (12)	0.0023 (12)
C123	0.0172 (13)	0.0185 (15)	0.0192 (14)	-0.0014 (11)	0.0031 (11)	0.0010 (11)
C124	0.0169 (11)	0.0207 (14)	0.0142 (11)	-0.0069 (11)	0.0033 (9)	-0.0006 (11)
C125	0.0194 (13)	0.0315 (16)	0.0168 (13)	-0.0005 (11)	0.0008 (10)	-0.0035 (12)

C126	0.0296 (15)	0.0371 (17)	0.0165 (13)	-0.0058 (13)	-0.0007 (11)	-0.0037 (12)
C127	0.0262 (14)	0.0346 (18)	0.0165 (13)	-0.0097 (13)	-0.0053 (11)	0.0073 (13)
C128	0.0244 (14)	0.0298 (16)	0.0255 (16)	-0.0001 (13)	-0.0022 (12)	0.0082 (13)
C129	0.0281 (15)	0.0206 (15)	0.0175 (14)	-0.0001 (12)	0.0003 (12)	0.0005 (11)
C130	0.0249 (14)	0.0187 (14)	0.0127 (12)	-0.0008 (11)	0.0042 (11)	-0.0021 (11)
C131	0.0208 (14)	0.0238 (16)	0.0269 (16)	0.0015 (12)	0.0003 (12)	-0.0045 (13)
C132	0.0346 (18)	0.034 (2)	0.0329 (19)	0.0137 (15)	0.0011 (15)	-0.0061 (15)
C133	0.052 (2)	0.0227 (16)	0.0290 (18)	0.0115 (15)	0.0079 (16)	0.0024 (14)
C134	0.055 (2)	0.0170 (14)	0.0257 (16)	-0.0024(14)	0.0108 (15)	-0.0011 (12)
C135	0.0345 (16)	0.0216 (14)	0.0214 (14)	-0.0052(12)	0.0074 (12)	-0.0013(12)
Ag4	0.02132 (10)	0.01748 (13)	0.01248 (10)	-0.00586 (9)	0.00067 (8)	0.00058 (8)
Cl4	0.0177 (3)	0.0165 (3)	0.0150 (3)	-0.0013(3)	-0.0017(2)	-0.0006(3)
N16	0.0125(11)	0.0154 (13)	0.0188(13)	-0.0024(9)	0.0018(10)	0.0033(10)
N17	0.0164 (11)	0.0153 (13)	0.0148 (12)	-0.0001(10)	0.0029 (10)	0.0010 (10)
N18	0.0154 (11)	0.0245 (14)	0.0176(12)	0.0005 (10)	0.0022(10)	0.0051 (11)
N19	0.0164(12)	0.0242(16)	0.0237(14)	-0.0041(11)	0.0022(10) 0.0024(11)	0.0001(11) 0.0047(12)
N20	0.0252(13)	0.0274(15)	0.0237(11) 0.0175(13)	-0.0065(11)	-0.0001(11)	0.0017(12) 0.0032(11)
P7	0.0252(15) 0.0158(3)	0.0271(13) 0.0142(4)	0.0173(13) 0.0147(3)	-0.0013(3)	0.0001(11) 0.0029(3)	0.0002(11)
P8	0.0133(4)	0.0179(4)	0.0114(3)	-0.0063(3)	0.0025(3)	-0.0004(3)
S4	0.00000(1)	0.0226(4)	0.0116(3)	-0.0037(3)	0.0024(3)	0.0026 (3)
C136	0.0156(12)	0.0220(1) 0.0177(14)	0.0130(3) 0.0148(13)	0.0000(11)	0.0021(3) 0.0023(11)	-0.0020(3)
C137	0.0165(12)	0.0138(14)	0.0167(14)	-0.0014(11)	0.0022(11)	-0.00020(11)
C138	0.0102(13)	0.0221(17)	0.0186(15)	-0.0034(12)	-0.0015(12)	0.0002(11) 0.0037(13)
C139	0.0172(13)	0.0221(17) 0.0167(14)	0.0137(13)	-0.0017(11)	0.0017(11)	-0.0029(11)
C140	0.0102(15) 0.0189(15)	0.0167(11)	0.0197(19) 0.0295(18)	-0.0103(15)	0.0017(11) 0.0058(14)	0.0029(11)
C141	0.0245(16)	0.050(3) 0.043(2)	0.0299(10) 0.0264(17)	-0.0117(15)	0.0036(11) 0.0076(14)	0.0044(15)
C142	0.0279(13)	0.0202(16)	0.0261(17) 0.0163(14)	-0.0042(12)	0.0012(11)	-0.0020(12)
C143	0.0184(14)	0.0202(10)	0.0289(18)	0.0080(16)	0.0012(11) 0.0013(13)	0.0262(12)
C144	0.054(3)	0.141(5)	0.0209(10)	0.068(3)	0.0013(2)	0.0202 (10) 0.009 (3)
C145	0.0284(15)	0.0180(15)	0.035(2)	-0.0018(11)	0.013(2) 0.0075(12)	-0.0017(11)
C146	0.0201(15) 0.0303(15)	0.0194 (14)	0.0280(15)	-0.0043(12)	0.0073(12)	-0.0035(12)
C147	0.0303(19) 0.0444(19)	0.0216(15)	0.0200(10) 0.0378(18)	-0.0027(14)	0.0225(16)	-0.0033(13)
C148	0.068 (2)	0.0283(18)	0.0233 (16)	-0.0047(17)	0.0242(17)	-0.0056(14)
C149	0.000(2) 0.054(2)	0.0205(18)	0.0255(15)	-0.0041(16)	0.0212(17) 0.0069(15)	-0.0040(13)
C150	0.0311(16)	0.0300(10) 0.0317(18)	0.0206(15)	-0.0050(13)	0.0034(13)	-0.0019(13)
C151	0.0183(12)	0.0213(15)	0.0189(14)	-0.0028(11)	0.0030(10)	0.0032(12)
C152	0.0218(13)	0.0204 (14)	0.0233 (14)	0.0030(11)	0.0023(11)	-0.0012(11)
C153	0.0237(15)	0.0201(17)	0.0253(11) 0.0273(16)	0.0052(13)	-0.0005(12)	0.0012(11) 0.0046(13)
C155	0.0237(13) 0.0232(14)	0.0369(18)	0.0264 (16)	-0.0040(13)	-0.0053(12)	0.0010(13)
C155	0.0252(11) 0.0368(17)	0.0284 (16)	0.0201(10) 0.0295(17)	-0.0066(14)	-0.0060(12)	-0.0018(13)
C156	0.0303(15)	0.0227 (15)	0.0239(15)	-0.0021(12)	-0.0052(12)	-0.0012(12)
C157	0.0162(12)	0.0227(10)	0.0237(13)	-0.0011(10)	0.0002(12)	-0.0022(11)
C158	0.0102(12) 0.0233(13)	0.0212(14)	0.0242(14)	0.0002(11)	0.00000(10)	0.0054(12)
C159	0.0255(15) 0.0251(14)	0.0212(11) 0.0199(14)	0.0266(15)	-0.0002(11)	0.0023(11)	0.0058(12)
C160	0.0293 (15)	0.0194 (14)	0.0206 (14)	0.0054 (12)	-0.0020(12)	0.0015(12)
C161	0.0254(14)	0.0298 (16)	0.0219 (15)	0.0047(12)	0.0025 (11)	-0.0027(12)
C162	0.0226 (13)	0.0210 (14)	0.0234 (15)	-0.0021(11)	0.0032 (11)	-0.0001(11)
C163	0.0274 (14)	0.0168 (14)	0.0130 (13)	0.0014 (11)	0.0065 (11)	-0.0003(11)
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C164	0.0319 (15)	0.0202 (15)	0.0146 (13)	0.0003 (12)	0.0065 (12)	-0.0005 (11)
C165	0.0299 (16)	0.0259 (17)	0.0193 (15)	0.0032 (13)	0.0047 (13)	0.0026 (12)
C166	0.0218 (14)	0.0315 (18)	0.0251 (16)	0.0043 (13)	0.0074 (13)	0.0025 (14)
C167	0.0271 (15)	0.0273 (16)	0.0196 (14)	0.0047 (13)	0.0079 (12)	0.0063 (13)
C168	0.0259 (15)	0.0194 (15)	0.0182 (14)	0.0022 (12)	0.0057 (12)	0.0044 (12)
C169	0.058 (2)	0.0185 (14)	0.0149 (14)	-0.0035 (15)	0.0044 (14)	0.0015 (12)
C170	0.077 (3)	0.0235 (16)	0.0295 (18)	-0.0127 (17)	-0.0135 (18)	0.0062 (14)
C171	0.134 (4)	0.0237 (18)	0.029 (2)	-0.019 (2)	-0.018 (2)	-0.0024 (15)
C172	0.133 (5)	0.0194 (18)	0.033 (2)	0.002 (2)	0.002 (3)	-0.0023 (16)
C173	0.091 (3)	0.0270 (19)	0.038 (2)	0.015 (2)	0.022 (2)	0.0006 (16)
C174	0.068 (3)	0.0170 (15)	0.0266 (17)	-0.0010 (16)	0.0137 (17)	0.0021 (13)
C175	0.0344 (15)	0.0246 (16)	0.0137 (12)	-0.0121 (14)	0.0004 (11)	0.0028 (12)
C176	0.0351 (17)	0.0351 (18)	0.0154 (14)	-0.0094 (14)	0.0037 (12)	-0.0015 (13)
C177	0.0383 (18)	0.0376 (18)	0.0202 (15)	-0.0051 (15)	0.0000 (13)	-0.0003 (13)
C178	0.0386 (18)	0.0322 (18)	0.0198 (15)	-0.0118 (14)	-0.0064 (13)	0.0040 (13)
C179	0.0510 (19)	0.0249 (17)	0.0105 (12)	-0.0118 (15)	-0.0027 (12)	-0.0007 (12)
C180	0.0490 (19)	0.0231 (15)	0.0182 (14)	-0.0037 (14)	0.0034 (13)	-0.0022 (12)

Geometric parameters (Å, °)

Ag1—P1	2.4727 (8)	Ag3—P5	2.4862 (8)
Ag1—P2	2.4877 (8)	Ag3—P6	2.4869 (8)
Ag1—Cl1	2.5905 (8)	Ag3—Cl3	2.5900 (8)
Ag1—S1	2.7192 (9)	Ag3—S3	2.6816 (8)
N1—N2	1.359 (3)	N11—C91	1.366 (4)
N1—C1	1.361 (4)	N11—N12	1.384 (4)
N1—H1N	0.875 (17)	N11—H11N	0.882 (17)
N2—C2	1.289 (4)	N12—C92	1.296 (4)
N3—C1	1.328 (4)	N13—C91	1.329 (4)
N3—C8	1.468 (4)	N13—C98	1.463 (4)
N3—H3N	0.871 (17)	N13—H13N	0.878 (17)
N4—C5	1.328 (4)	N14—C94	1.338 (4)
N4—C4	1.345 (4)	N14—C95	1.348 (4)
N5—C7	1.325 (4)	N15—C97	1.335 (4)
N5—C6	1.339 (4)	N15—C96	1.340 (4)
P1-C16	1.821 (3)	P5—C106	1.824 (3)
P1-C10	1.835 (3)	P5—C112	1.825 (3)
P1—C22	1.838 (3)	P5-C100	1.833 (3)
P2—C34	1.822 (3)	P6—C130	1.829 (3)
P2—C28	1.824 (3)	P6—C118	1.830 (3)
P2—C40	1.842 (3)	P6—C124	1.831 (3)
S1—C1	1.707 (3)	S3—C91	1.700 (3)
C2—C4	1.483 (4)	С92—С94	1.486 (4)
C2—C3	1.501 (5)	С92—С93	1.493 (5)
С3—НЗА	0.9800	С93—Н93А	0.9800
С3—Н3В	0.9800	С93—Н93В	0.9800
С3—НЗС	0.9800	С93—Н93С	0.9800
C4—C7	1.400 (4)	C94—C97	1.406 (4)

C5—C6	1.391 (5)	C95—C96	1.373 (5)
С5—Н5	0.9500	С95—Н95	0.9500
С6—Н6	0.9500	С96—Н96	0.9500
С7—Н7	0.9500	С97—Н97	0.9500
C8—C9	1.494 (5)	С98—С99	1.512 (5)
C8—H8A	0.9900	С98—Н98А	0.9900
C8—H8B	0.9900	С98—Н98В	0.9900
С9—Н9А	0.9800	С99—Н99А	0.9800
С9—Н9В	0.9800	С99—Н99В	0.9800
С9—Н9С	0.9800	С99—Н99С	0.9800
C10—C15	1.389 (4)	C100—C105	1.392 (4)
C10—C11	1.398 (4)	C100—C101	1.405 (4)
C11—C12	1.399 (4)	C101—C102	1.391 (4)
C11—H11	0.9500	C101—H101	0.9500
C12—C13	1.377 (4)	C102—C103	1.388 (4)
C12—H12	0.9500	C102—H102	0.9500
C13—C14	1.386 (4)	C103—C104	1.387 (5)
С13—Н13	0.9500	C103—H103	0.9500
C14—C15	1.393 (4)	C104—C105	1.392 (4)
C14—H14	0.9500	C104—H104	0.9500
С15—Н15	0.9500	C105—H105	0.9500
C16—C21	1.389 (4)	C106—C107	1.393 (4)
C16—C17	1.394 (4)	C106—C111	1.403 (4)
C17—C18	1.389 (4)	C107—C108	1.392 (4)
С17—Н17	0.9500	C107—H107	0.9500
C18—C19	1.393 (5)	C108—C109	1.383 (4)
C18—H18	0.9500	C108—H108	0.9500
C19—C20	1.373 (5)	C109—C110	1.401 (4)
С19—Н19	0.9500	C109—H109	0.9500
C20—C21	1.396 (5)	C110—C111	1.392 (4)
С20—Н20	0.9500	C110—H110	0.9500
C21—H21	0.9500	C111—H111	0.9500
C22—C27	1.390 (4)	C112—C117	1.392 (4)
C22—C23	1.396 (4)	C112—C113	1.395 (4)
C23—C24	1.400 (4)	C113—C114	1.401 (4)
С23—Н23	0.9500	C113—H113	0.9500
C24—C25	1.379 (5)	C114—C115	1.379 (4)
C24—H24	0.9500	C114—H114	0.9500
C25—C26	1.377 (5)	C115—C116	1.377 (4)
С25—Н25	0.9500	C115—H115	0.9500
C26—C27	1.396 (4)	C116—C117	1.392 (4)
С26—Н26	0.9500	C116—H116	0.9500
С27—Н27	0.9500	C117—H117	0.9500
C28—C29	1.399 (4)	C118—C123	1.391 (4)
C28—C33	1.404 (5)	C118—C119	1.395 (4)
C29—C30	1.394 (4)	C119—C120	1.398 (4)
С29—Н29	0.9500	С119—Н119	0.9500
C30—C31	1.381 (5)	C120—C121	1.392 (5)

С30—Н30	0.9500	C120—H120	0.9500
C31—C32	1.384 (5)	C121—C122	1.390 (5)
C31—H31	0.9500	C121—H121	0.9500
C32—C33	1.393 (5)	C122—C123	1.397 (4)
С32—Н32	0.9500	C122—H122	0.9500
С33—Н33	0.9500	C123—H123	0.9500
C34—C35	1.385 (4)	C124—C125	1.383 (4)
C34—C39	1,409 (4)	C124—C129	1.388 (4)
C35—C36	1.393 (5)	C125—C126	1.388 (3)
С35—Н35	0.9500	C125—H125	0.9500
$C_{36} - C_{37}$	1 371 (4)	C126—C127	1 382 (4)
C36—H36	0.9500	C126—H126	0.9500
C_{37} C_{38}	1 387 (5)	$C_{120} = C_{120}$	1.379(4)
C37—E38	0.9500	C127—C128	0.9500
C_{3}^{2} C_{3}^{2}	1.386(A)	$C_{12} = C_{12}$	1.305(4)
C_{20} H_{20}	0.0500	$C_{120} = C_{120}$	1.393 (4)
C30_H30	0.9500	С120—1120	0.9300
C39—H39	0.9500	C129—H129	0.9500
	1.383 (5)	C130—C131	1.385 (4)
C40—C45	1.395 (5)		1.400 (4)
C41—C42	1.404 (5)	C131—C132	1.392 (5)
C41—H41	0.9500	С131—Н131	0.9500
C42—C43	1.363 (6)	C132—C133	1.383 (5)
C42—H42	0.9500	С132—Н132	0.9500
C43—C44	1.398 (6)	C133—C134	1.373 (5)
C43—H43	0.9500	С133—Н133	0.9500
C44—C45	1.396 (5)	C134—C135	1.391 (4)
C44—H44	0.9500	C134—H134	0.9500
C45—H45	0.9500	С135—Н135	0.9500
Ag2—P3	2.4858 (8)	Ag4—P7	2.4784 (8)
Ag2—P4	2.4861 (8)	Ag4—P8	2.4838 (8)
Ag2—Cl2	2.5932 (8)	Ag4—Cl4	2.5901 (8)
Ag2—S2	2.6848 (8)	Ag4—S4	2.7412 (9)
N6—N7	1.362 (3)	N16—C136	1.364 (4)
N6—C46	1.363 (4)	N16—N17	1.387 (3)
N6—H6N	0.877 (17)	N16—H16N	0.860 (17)
N7—C47	1.287 (4)	N17—C137	1.300 (4)
N8—C46	1.330 (4)	N18—C136	1.325 (4)
N8—C53	1.465 (4)	N18—C143	1.456 (4)
N8—H8N	0.861 (17)	N18—H18N	0.860 (17)
N9—C50	1.337 (4)	N19—C139	1.339 (4)
N9—C49	1.347 (4)	N19—C140	1.343 (4)
N10-C52	1.328 (4)	N20-C142	1.336 (4)
N10-C51	1 343 (4)	N20-C141	1 344 (4)
P3—C61	1.833 (3)	P7—C145	1.823 (3)
P3—C55	1 834 (3)	P7—C157	1 826 (3)
P3C67	1 838 (3)	P7—C151	1 832 (3)
P4—C85	1 822 (3)	P8—C163	1 825 (3)
P4C73	1 824 (3)	P8-C169	1.025(3) 1.825(3)
11 015	1.047 (3)	10 0107	1.025 (5)

P4—C79	1.835 (3)	P8—C175	1.840 (3)
S2—C46	1.703 (3)	S4—C136	1.705 (3)
C47—C49	1.482 (4)	C137—C139	1.482 (4)
C47—C48	1.503 (4)	C137—C138	1.494 (4)
C48—H48A	0.9800	C138—H13A	0.9800
C48—H48B	0.9800	C138—H13B	0.9800
C48—H48C	0.9800	C138—H13C	0.9800
C49-C52	1 404 (4)	C139-C142	1410(5)
C_{50} C_{51}	1 386 (5)	C140-C141	1.110(5)
C50—H50	0.9500	C_{140} H_{140}	0.9500
C51 H51	0.9500	C_{141} H141	0.9500
C52 H52	0.9500	$C_{141} = 1141$	0.9500
C52—H52	0.9300	C142	1.524 (()
C53-C54	1.494 (4)	C143 - C144	1.324 (0)
C53—H53A	0.9900	С143—П14А	0.9900
С53—Н53В	0.9900	C143—H14B	0.9900
С54—Н54А	0.9800	C144—H14C	0.9800
C54—H54B	0.9800	C144—H14D	0.9800
C54—H54C	0.9800	C144—H14E	0.9800
C55—C56	1.383 (4)	C145—C146	1.393 (4)
C55—C60	1.396 (4)	C145—C150	1.403 (5)
C56—C57	1.382 (4)	C146—C147	1.395 (4)
С56—Н56	0.9500	C146—H146	0.9500
C57—C58	1.375 (4)	C147—C148	1.385 (5)
С57—Н57	0.9500	C147—H147	0.9500
C58—C59	1.398 (4)	C148—C149	1.392 (5)
C58—H58	0.9500	C148—H148	0.9500
C59—C60	1.389 (4)	C149—C150	1.389 (5)
С59—Н59	0.9500	C149—H149	0.9500
С60—Н60	0.9500	C150—H150	0.9500
C61—C66	1.389 (4)	C151—C152	1.396 (4)
C61—C62	1.397 (4)	C151—C156	1.398 (4)
C62—C63	1.397 (4)	C152—C153	1.397 (4)
C62 - H62	0.9500	C152—H152	0.9500
C63 - C64	1 396 (4)	C_{153} C_{154}	1.380(5)
C63_H63	0.9500	C153—H153	0.9500
C64 C65	1.371(4)	C154 C155	1 301 (5)
C64 H64	0.0500	$C_{154} = C_{155}$	0.0500
C65 C66	0.9500	$C_{154} = 11154$	1.204(4)
C(5) = C(6)	1.393 (4)	C155_U155	1.394 (4)
C65—H65	0.9500	С155—Н155	0.9500
C66—H66	0.9500	С156—Н156	0.9500
C67—C72	1.387 (4)	C157—C162	1.393 (4)
C67—C68	1.393 (4)	C157—C158	1.404 (4)
C68—C69	1.399 (4)	C158—C159	1.397 (4)
С68—Н68	0.9500	C158—H158	0.9500
C69—C70	1.389 (4)	C159—C160	1.391 (4)
С69—Н69	0.9500	C159—H159	0.9500
C70—C71	1.380 (5)	C160—C161	1.381 (4)
С70—Н70	0.9500	C160—H160	0.9500

C71—C72	1.393 (4)	C161—C162	1.397 (4)
C71—H71	0.9500	C161—H161	0.9500
С72—Н72	0.9500	C162—H162	0.9500
C73—C74	1.401 (4)	C163—C164	1.393 (4)
C73—C78	1.401 (4)	C163—C168	1.394 (4)
C74—C75	1.392 (4)	C164—C165	1.392 (5)
С74—Н74	0.9500	C164—H164	0.9500
C75—C76	1.384 (5)	C165—C166	1.387 (5)
С75—Н75	0.9500	C165—H165	0.9500
C76—C77	1.385 (5)	C166—C167	1.386 (5)
С76—Н76	0.9500	C166—H166	0.9500
С77—С78	1.392 (4)	C167—C168	1.395 (4)
С77—Н77	0.9500	С167—Н167	0.9500
С78—Н78	0.9500	C168—H168	0.9500
С79—С80	1.391 (4)	C169—C174	1.392 (5)
С79—С84	1.399 (4)	C169—C170	1.407 (5)
C80—C81	1.405 (5)	C170—C171	1.407 (5)
С80—Н80	0.9500	С170—Н170	0.9500
C81—C82	1.374 (5)	C171—C172	1.396 (7)
C81—H81	0.9500	C171—H171	0.9500
C82—C83	1.386 (5)	C172—C173	1.369 (7)
C82—H82	0.9500	С172—Н172	0.9500
C83—C84	1.388 (4)	C173—C174	1.405 (5)
С83—Н83	0.9500	С173—Н173	0.9500
C84—H84	0.9500	C174—H174	0.9500
C85—C90	1.397 (4)	C175—C176	1.380 (5)
C85—C86	1.400 (4)	C175—C180	1.401 (4)
C86—C87	1.391 (4)	C176—C177	1.409 (5)
C86—H86	0.9500	С176—Н176	0.9500
C87—C88	1.373 (5)	C177—C178	1.378 (5)
C87—H87	0.9500	С177—Н177	0.9500
C88—C89	1.379 (5)	C178—C179	1.369 (5)
C88—H88	0.9500	C178—H178	0.9500
C89—C90	1.391 (4)	C179—C180	1.394 (4)
C89—H89	0.9500	C179—H179	0.9500
C90—H90	0.9500	C180—H180	0.9500
	0.9200		0.9000
P1—Ag1—P2	115.46 (3)	P5—Ag3—P6	118.91 (3)
P1 - Ag1 - C11	119.77 (3)	P5 - Ag3 - Cl3	119.36 (3)
P2—Ag1—C11	116.02 (3)	P6—Ag3—Cl3	112.62 (2)
P1 - Ag1 - S1	99.69 (3)	P5—Ag3—S3	97.41 (2)
P2—Ag1—S1	112.46 (3)	P6—Ag3—S3	112.54 (3)
Cl1 - Ag1 - S1	87.76 (3)	Cl3—Ag3—S3	90.32 (3)
N2—N1—C1	119.0 (3)	C91—N11—N12	117.6 (3)
N2—N1—H1N	121 (2)	C91—N11—H11N	118 (2)
C1—N1—H1N	114 (2)	N12—N11—H11N	125 (2)
C2—N2—N1	117.7 (3)	C92—N12—N11	116.9 (3)
C1—N3—C8	123.4 (3)	C91—N13—C98	123.1 (3)

C1—N3—H3N	117 (2)	C91—N13—H13N	122 (2)
C8—N3—H3N	119 (2)	C98—N13—H13N	115 (2)
C5—N4—C4	116.6 (3)	C94—N14—C95	116.0 (3)
C7—N5—C6	116.3 (3)	C97—N15—C96	116.0 (3)
C16—P1—C10	102.86 (13)	C106—P5—C112	103.53 (14)
C16—P1—C22	104.30 (13)	C106—P5—C100	101.08 (13)
C10—P1—C22	103.17 (14)	C112—P5—C100	106.26 (13)
C16—P1—Ag1	113.27 (10)	C106—P5—Ag3	119.87 (9)
C10—P1—Ag1	121.53 (10)	C112—P5—Ag3	106.84 (10)
C22—P1—Ag1	110.00 (10)	C100—P5—Ag3	117.75 (10)
C34—P2—C28	106.60 (14)	C130—P6—C118	103.54 (14)
C_{34} P2 C40	103.49 (15)	C130—P6—C124	102.20 (14)
$C_{28} = P_{2} = C_{40}$	102.28 (15)	$C_{118} - P_{6} - C_{124}$	103.33(13)
C_{34} P2 Ag1	112.92 (10)	C130 P6 Ag3	113 15 (9)
C_{28} P2 Ag1	116.85 (11)	C118—P6—Ag3	117.96 (10)
C40 - P2 - Ag1	113 34 (10)	C124 P6 Ag3	114 79 (9)
$C_1 = S_1 = A_{g_1}$	99.81 (10)	$C_{124} = 10^{-1} Mg^{3}$	102.46(10)
$N_{1} = N_{1}$	117 3 (3)	N13 C01 N11	102.40(10) 117.4(3)
$N_2 = C_1 = N_1$	117.5(3) 123.6(2)	N13 C01 S3	117.4(3) 122.3(2)
$N_1 = C_1 = S_1$	123.0(2)	N11 C01 S2	123.3(2)
N2 C2 C4	119.2(2) 115.1(2)	N12 = C02 = C04	119.3(2) 114.3(3)
$N_2 = C_2 = C_4$	113.1(3) 124.2(2)	N12 - C92 - C94	114.3(3)
$N_2 = C_2 = C_3$	124.5 (3)	N12 - C92 - C93	125.5(3)
C4-C2-C3	120.6 (3)	C94 - C92 - C93	120.4 (3)
C2—C3—H3A	109.5	С92—С93—Н93А	109.5
C2—C3—H3B	109.5	С92—С93—Н93В	109.5
НЗА—СЗ—НЗВ	109.5	Н93А—С93—Н93В	109.5
C2—C3—H3C	109.5	С92—С93—Н93С	109.5
НЗА—СЗ—НЗС	109.5	Н93А—С93—Н93С	109.5
НЗВ—СЗ—НЗС	109.5	Н93В—С93—Н93С	109.5
N4—C4—C7	120.2 (3)	N14—C94—C97	121.0 (3)
N4—C4—C2	118.3 (3)	N14—C94—C92	117.8 (3)
C7—C4—C2	121.5 (3)	C97—C94—C92	121.3 (3)
N4—C5—C6	122.7 (3)	N14—C95—C96	122.5 (3)
N4—C5—H5	118.7	N14—C95—H95	118.7
С6—С5—Н5	118.7	С96—С95—Н95	118.7
N5—C6—C5	121.0 (3)	N15—C96—C95	122.0 (3)
N5—C6—H6	119.5	N15—C96—H96	119.0
С5—С6—Н6	119.5	С95—С96—Н96	119.0
N5—C7—C4	123.0 (3)	N15—C97—C94	122.3 (3)
N5—C7—H7	118.5	N15—C97—H97	118.9
С4—С7—Н7	118.5	С94—С97—Н97	118.9
N3—C8—C9	113.0 (3)	N13—C98—C99	111.9 (3)
N3—C8—H8A	109.0	N13—C98—H98A	109.2
С9—С8—Н8А	109.0	С99—С98—Н98А	109.2
N3—C8—H8B	109.0	N13—C98—H98B	109.2
С9—С8—Н8В	109.0	С99—С98—Н98В	109.2
H8A—C8—H8B	107.8	H98A—C98—H98B	107.9
С8—С9—Н9А	109.5	С98—С99—Н99А	109.5

С8—С9—Н9В	109.5	С98—С99—Н99В	109.5
Н9А—С9—Н9В	109.5	Н99А—С99—Н99В	109.5
С8—С9—Н9С	109.5	С98—С99—Н99С	109.5
H9A—C9—H9C	109.5	Н99А—С99—Н99С	109.5
H9B—C9—H9C	109.5	Н99В—С99—Н99С	109.5
C15—C10—C11	120.1 (3)	C105—C100—C101	119.1 (3)
C15—C10—P1	121.5 (2)	C105—C100—P5	118.4 (2)
C11—C10—P1	118.3 (2)	C101—C100—P5	122.4 (2)
C10—C11—C12	119.5 (3)	C102—C101—C100	120.9 (3)
C10—C11—H11	120.2	C102—C101—H101	119.5
C12—C11—H11	120.2	C100-C101-H101	119.5
C13—C12—C11	119.9 (3)	C103—C102—C101	119.4 (3)
C13—C12—H12	120.0	C103—C102—H102	120.3
C11—C12—H12	120.0	C101—C102—H102	120.3
C12—C13—C14	120.7 (3)	C104—C103—C102	119.9 (3)
С12—С13—Н13	119.6	C104—C103—H103	120.1
C14—C13—H13	119.6	C102 - C103 - H103	120.1
C13 - C14 - C15	119.9 (3)	C102 - C103 - C104 - C105	120.1 1211(3)
C_{13} C_{14} H_{14}	120.1	C103 - C104 - H104	119.4
C_{15} C_{14} H_{14}	120.1	C105 - C104 - H104	119.1
C10-C15-C14	119.8 (3)	C100-C105-C104	119.5 (3)
C10-C15-H15	120.1	C100 - C105 - H105	120.3
C_{14} C_{15} H_{15}	120.1	C104 - C105 - H105	120.3
C_{21} C_{16} C_{17}	118.9 (3)	C107 - C106 - C111	120.3 110.2(3)
$C_{21} = C_{10} = C_{17}$	1177(2)	C107 - C106 - P5	117.2(3) 117.8(2)
C_{17} C_{16} P_{1}	117.7(2) 123 4 (2)	$C_{111} - C_{106} - P_5$	117.0(2) 122.9(2)
C_{18} C_{17} C_{16}	123.7(2) 120.3(3)	$C_{108} = C_{107} = C_{106}$	122.9(2) 120.5(3)
$C_{18} = C_{17} = C_{10}$	110.0	C108 - C107 - C100	120.3 (3)
$C_{16} = C_{17} = H_{17}$	119.9	C106 - C107 - H107	119.7
$C_{17} = C_{17} = C$	120.1 (3)	C100 - C108 - C107	119.7 120.3(3)
$C_{17} = C_{18} = C_{19}$	120.1 (5)	C109 - C108 - C107	110.8
C19 - C18 - H18	120.0	C107 - C108 - H108	119.0
C_{20} C_{19} C_{18}	120.0 (3)	C108 - C109 - C110	119.6(3)
$C_{20} = C_{19} = C_{18}$	120.0 (3)	C108 - C109 - C110	119.0 (3)
$C_{20} = C_{19} = H_{19}$	120.0	$C_{100} = C_{100} = H_{100}$	120.2
$C_{10} = C_{10} = C_{11}$	120.0 110.0(A)	$C_{111} = C_{110} = C_{109}$	120.2 120.2(3)
$C_{19} = C_{20} = C_{21}$	120.0		110.0
$C_{1}^{2} = C_{2}^{2} = C_{12}^{2} = C_{2}^{2} = C_{$	120.0	$C_{100} = C_{110} = H_{110}$	119.9
$C_{21} = C_{20} = 1120$	120.0 120.7(2)	$C_{10} = C_{110} = C_{106}$	117.7 120.0(2)
$C_{10} = C_{21} = C_{20}$	120.7 (3)	$C_{110} = C_{111} = C_{100}$	120.0 (3)
$C_{10} = C_{21} = H_{21}$	119.0	C106 C111 H111	120.0
$C_{20} = C_{21} = H_{21}$	119.0	$C_{110} - C_{111} - C_{112}$	120.0 118.2(3)
$C_{27} = C_{22} = C_{23}$	118.0(3)	$C_{117} = C_{112} = C_{113}$	116.2(3)
$C_{27} = C_{22} = 11$	110.0(2) 122.1(2)	$C_{117} - C_{112} - C_{15}$	110.0(2) 125.2(2)
$C_{23} = C_{22} = C_{13}$	123.1(2) 120.0(2)	$C_{113} - C_{112} - C_{114}$	123.2(2) 120.5(2)
$\begin{array}{cccccccccccccccccccccccccccccccccccc$	120.0 (5)	$C_{112} = C_{113} = C_{114}$	120.3 (3)
$C_{22} = C_{23} = \Pi_{23}$	120.0	С112—С113—П113	119.0
$C_{24} = C_{23} = H_{23}$	120.0		119.8
$U_{23} - U_{24} - U_{23}$	120.0 (3)	UII3—UII4—UII3	120.0(3)

C25—C24—H24	119.7	C115—C114—H114	120.0
C23—C24—H24	119.7	C113—C114—H114	120.0
C26—C25—C24	119.5 (3)	C116—C115—C114	120.3 (3)
С26—С25—Н25	120.2	C116—C115—H115	119.9
C24—C25—H25	120.2	C114—C115—H115	119.9
C_{25} C_{26} C_{27}	120.5(3)	$C_{115} - C_{116} - C_{117}$	119.8 (3)
$C_{25} = C_{26} = H_{26}$	119.8	C115—C116—H116	120.1
C_{27} C_{26} H_{26}	119.8	C117_C116_H116	120.1
C_{22}^{22} C_{23}^{22} C_{25}^{22}	120 5 (3)	C_{112} C_{117} C_{116}	120.1 121.2(3)
$C_{22} = C_{27} = C_{20}$	110.8	$C_{112} = C_{117} = C_{110}$	121.2 (3)
$C_{22} = C_{27} = H_{27}$	119.8	$C_{112} - C_{117} - H_{117}$	119.4
$C_{20} = C_{27} = H_{27}$	117.0 119.7(2)	$C_{110} - C_{117} - III17$	119.4
$C_{29} = C_{28} = C_{33}$	110.7 (3)	C123—C118—C119	119.7 (3)
C29—C28—P2	123.3 (3)	C123—C118—P6	122.5 (2)
C33—C28—P2	118.0 (2)	C119—C118—P6	117.8(2)
C30—C29—C28	120.5 (3)	C118—C119—C120	119.9 (3)
С30—С29—Н29	119.8	C118—C119—H119	120.0
С28—С29—Н29	119.8	C120—C119—H119	120.0
C31—C30—C29	120.1 (3)	C121—C120—C119	120.3 (3)
С31—С30—Н30	120.0	C121—C120—H120	119.9
С29—С30—Н30	120.0	C119—C120—H120	119.9
C30—C31—C32	120.2 (3)	C122—C121—C120	119.7 (3)
С30—С31—Н31	119.9	C122—C121—H121	120.2
C32—C31—H31	119.9	C120—C121—H121	120.2
C31—C32—C33	120.3 (3)	C121—C122—C123	120.2 (3)
С31—С32—Н32	119.9	C121—C122—H122	119.9
С33—С32—Н32	119.9	C123—C122—H122	119.9
$C_{32} - C_{33} - C_{28}$	120.2 (3)	C118 - C123 - C122	120.2 (3)
C32—C33—H33	119.9	$C_{118} - C_{123} - H_{123}$	119.9
C28—C33—H33	119.9	$C_{122} - C_{123} - H_{123}$	119.9
C_{35} C_{34} C_{39}	117.6 (3)	C_{125} C_{124} C_{129}	119.1 (3)
$C_{35} = C_{34} = C_{35}$	117.0(3) 118.2(2)	$C_{125} = C_{124} = C_{125}$	117.1(3) 122.7(2)
$C_{39} = C_{34} = P_{2}$	110.2(2) 124.2(2)	$C_{123} - C_{124} - P_6$	122.7(2) 118.2(2)
$C_{3} = C_{3} = C_{3} = C_{3}$	124.2(2) 121.5(2)	$C_{12}^{12} = C_{12}^{12} = C_{12}^{10}$	110.2(2)
$C_{24} = C_{25} = C_{30}$	121.3 (3)	C124 - C125 - C120	120.4 (3)
С34—С35—П35	119.5	C124—C125—H125	119.8
C36—C35—H35	119.3	C126—C125—H125	119.8
$C_{37} - C_{36} - C_{35}$	120.1 (3)	C12/-C126-C125	120.36 (16)
С37—С36—Н36	119.9	C127—C126—H126	119.8
С35—С36—Н36	119.9	C125—C126—H126	119.8
C36—C37—C38	119.8 (3)	C128—C127—C126	119.70 (17)
С36—С37—Н37	120.1	C128—C127—H127	120.1
С38—С37—Н37	120.1	C126—C127—H127	120.1
C39—C38—C37	120.2 (3)	C127—C128—C129	120.0 (3)
С39—С38—Н38	119.9	C127—C128—H128	120.0
С37—С38—Н38	119.9	C129—C128—H128	120.0
C38—C39—C34	120.7 (3)	C124—C129—C128	120.4 (3)
С38—С39—Н39	119.6	C124—C129—H129	119.8
С34—С39—Н39	119.6	C128—C129—H129	119.8
C41—C40—C45	119.1 (3)	C131—C130—C135	117.8 (3)

C41—C40—P2	120.1 (3)	C131—C130—P6	124.1 (2)
C45—C40—P2	120.5 (3)	C135—C130—P6	118.1 (2)
C40—C41—C42	120.5 (4)	C130—C131—C132	121.2 (3)
C40—C41—H41	119.7	C130—C131—H131	119.4
C42—C41—H41	119.7	C132—C131—H131	119.4
C43—C42—C41	120.5 (4)	C133 - C132 - C131	119.8 (3)
C43-C42-H42	119.8	C133—C132—H132	120.1
C41—C42—H42	119.8	C131—C132—H132	120.1
C42-C43-C44	119.7 (4)	C_{134} C_{133} C_{132}	120.2(3)
C42 - C43 - H43	120.2	C134—C133—H133	119.9
C44-C43-H43	120.2	C132—C133—H133	119.9
$C_{45} - C_{44} - C_{43}$	120.2 120.1 (4)	C_{133} C_{134} C_{135}	119.8 (3)
$C_{45} = C_{44} = C_{45}$	110.0	$C_{133} = C_{134} = C_{135}$	120.1
$C_{43} = C_{44} = H_{44}$	119.9	$C_{135} = C_{134} = H_{134}$	120.1
$C_{43} = C_{44} = \Pi_{44}$	119.9 120.1 (4)	$C_{133} = C_{134} = C_{134}$	120.1 121.2(2)
C40 - C45 - U45	120.1 (4)	$C_{124} = C_{125} = C_{130}$	121.2(3)
C40 - C43 - H45	120.0	С134—С135—П135	119.4
C44 - C45 - H45	120.0	C130—C135—H135	119.4
P3—Ag2—P4	119.09 (3)	P/-Ag4-P8	113.84 (3)
P3—Ag2—Cl2	119.20 (3)	P/—Ag4—Cl4	118.99 (3)
P4—Ag2—Cl2	112.70 (2)	P8—Ag4—Cl4	117.37 (3)
P3—Ag2—S2	96.42 (3)	P7—Ag4—S4	102.70 (2)
P4—Ag2—S2	113.81 (3)	P8—Ag4—S4	111.64 (3)
Cl2—Ag2—S2	89.81 (3)	Cl4—Ag4—S4	87.39 (2)
N7—N6—C46	118.8 (3)	C136—N16—N17	116.9 (3)
N7—N6—H6N	120 (2)	C136—N16—H16N	118 (2)
C46—N6—H6N	115 (2)	N17—N16—H16N	125 (2)
C47—N7—N6	118.0 (3)	C137—N17—N16	116.7 (3)
C46—N8—C53	123.5 (3)	C136—N18—C143	124.1 (3)
C46—N8—H8N	118 (2)	C136—N18—H18N	121 (2)
C53—N8—H8N	119 (2)	C143—N18—H18N	114 (2)
C50—N9—C49	116.2 (3)	C139—N19—C140	115.6 (3)
C52—N10—C51	116.0 (3)	C142—N20—C141	115.0 (3)
C61—P3—C55	103.27 (14)	C145—P7—C157	102.91 (14)
C61—P3—C67	101.48 (13)	C145—P7—C151	104.25 (14)
C55—P3—C67	106.38 (13)	C157—P7—C151	103.01 (14)
C61—P3—Ag2	120.58 (9)	C145—P7—Ag4	112.01 (11)
C55—P3—Ag2	107.03 (10)	C157—P7—Ag4	123.37 (10)
C67 - P3 - Ag2	116.64 (10)	C151—P7—Ag4	109.43 (10)
C85 - P4 - C73	103 60 (13)	C163 - P8 - C169	102.70 (15)
C85—P4—C79	102.28 (14)	C163 - P8 - C175	106.46 (14)
C73 - P4 - C79	103.66 (14)	$C_{169} = P_{8} = C_{175}$	103 47 (16)
C85 P4 Ag2	114 03 (9)	$C_{163} = P_8 = A_{94}$	117.03(10)
C73 P4 Ag2	118.23(10)	C_{169} P8 Ag4	112 24 (10)
$C79 P4 A \sigma^2$	113 24 (10)	C175—P8—A g4	113 55 (10)
$C_{46} = S_{2} = \Delta \sigma^{2}$	102 38 (10)	$C136 S4 \Delta \sigma 4$	98 98 (10)
N8 - C46 - N6	117 1 (3)	$130 - 37 - Ag^{-1}$	117 4 (3)
N8 - C46 - S2	122 0 (2)	N18 - C136 - S4	127.7(3) 1240(2)
N6-C46-S2	122.9(2) 120.0(2)	N16 C136 S4	127.0(2) 1186(2)
110 010 52	120.0 (2)	110 0130 - S T	110.0 (2)

N7—C47—C49	115.3 (3)	N17—C137—C139	114.1 (3)
N7—C47—C48	124.5 (3)	N17—C137—C138	125.3 (3)
C49—C47—C48	120.2 (3)	C139—C137—C138	120.5 (3)
C47—C48—H48A	109.5	C137—C138—H13A	109.5
C47—C48—H48B	109.5	C137—C138—H13B	109.5
H48A—C48—H48B	109.5	H13A—C138—H13B	109.5
C47—C48—H48C	109.5	C137—C138—H13C	109.5
H48A—C48—H48C	109.5	H13A—C138—H13C	109.5
H48B—C48—H48C	109.5	H13B—C138—H13C	109.5
N9—C49—C52	120.5 (3)	N19—C139—C142	121.3 (3)
N9—C49—C47	118.1 (3)	N19—C139—C137	117.8 (3)
C52—C49—C47	121.3 (3)	C142—C139—C137	121.0 (3)
N9-C50-C51	122.6 (3)	N19-C140-C141	122.6(3)
N9-C50-H50	118 7	N19—C140—H140	118 7
$C_{51} - C_{50} - H_{50}$	118.7	C_{141} C_{140} H_{140}	118.7
N10-C51-C50	121.6 (3)	N20-C141-C140	122.9 (3)
N10-C51-H51	119.2	N20-C141-H141	118.6
C_{50} C_{51} H_{51}	119.2	C_{140} C_{141} H_{141}	118.6
N10 C52 C49	122 0 (3)	N20 C142 C130	122.3(3)
N10 - C52 - C49	122.9 (3)	$N_{20} = C_{142} = C_{153}$	122.3 (3)
$C_{10} = C_{52} = H_{52}$	118.6	120 - C142 - 11142	110.0
$N_{2}^{8} C_{3}^{53} C_{54}^{54}$	112 4 (2)	N18 C142 C144	100.4(3)
N9 C52 H52A	113.4 (3)	N18 - C143 - C144	109.4 (3)
No-C33-H33A	108.9	$N16 - C143 - \Pi14A$	109.8
C54—C53—H53A	108.9	C144—C145—H14A	109.8
N8-C33-H33B	108.9	N18—C143—H14B	109.8
С54—С53—Н53В	108.9	C144—C143—H14B	109.8
Н53А—С53—Н53В	107.7	H14A—C143—H14B	108.2
С53—С54—Н54А	109.5	C143— $C144$ — $H14C$	109.5
С53—С54—Н54В	109.5	C143—C144—H14D	109.5
H54A—C54—H54B	109.5	H14C—C144—H14D	109.5
С53—С54—Н54С	109.5	C143—C144—H14E	109.5
H54A—C54—H54C	109.5	H14C—C144—H14E	109.5
H54B—C54—H54C	109.5	H14D—C144—H14E	109.5
C56—C55—C60	118.8 (3)	C146—C145—C150	119.5 (3)
C56—C55—P3	116.6 (2)	C146—C145—P7	123.5 (2)
C60—C55—P3	124.6 (2)	C150—C145—P7	117.0 (2)
C57—C56—C55	121.2 (3)	C145—C146—C147	120.3 (3)
С57—С56—Н56	119.4	C145—C146—H146	119.8
С55—С56—Н56	119.4	C147—C146—H146	119.8
C58—C57—C56	120.3 (3)	C148—C147—C146	119.6 (3)
С58—С57—Н57	119.8	C148—C147—H147	120.2
С56—С57—Н57	119.8	C146—C147—H147	120.2
C57—C58—C59	119.4 (3)	C147—C148—C149	120.7 (3)
С57—С58—Н58	120.3	C147—C148—H148	119.7
С59—С58—Н58	120.3	C149—C148—H148	119.7
C60—C59—C58	120.2 (3)	C150—C149—C148	119.8 (4)
С60—С59—Н59	119.9	C150—C149—H149	120.1
С58—С59—Н59	119.9	C148—C149—H149	120.1

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C59—C60—C55	120.1 (3)	C149—C150—C145	120.0 (3)
С59—С60—Н60	119.9	C149—C150—H150	120.0
С55—С60—Н60	119.9	C145—C150—H150	120.0
C66—C61—C62	119.4 (3)	C152—C151—C156	118.0 (3)
C66—C61—P3	117.7 (2)	C152—C151—P7	118.3 (2)
C62—C61—P3	122.9 (2)	C156—C151—P7	123.7 (2)
C61—C62—C63	120.0 (3)	C151—C152—C153	120.9 (3)
С61—С62—Н62	120.0	C151—C152—H152	119.6
C63—C62—H62	120.0	C153—C152—H152	119.6
C64—C63—C62	119.9 (3)	C154—C153—C152	120.4(3)
C64 - C63 - H63	120.0	C154—C153—H153	119.8
C62 - C63 - H63	120.0	$C_{152} = C_{153} = H_{153}$	119.8
C65 - C64 - C63	119.8 (3)	C152 = C155 = T1155	119.0 119.5(3)
C65 C64 H64	119.8 (3)	C153 C154 H154	119.3 (5)
C63 $C64$ $H64$	120.1	C155 C154 H154	120.3
C64 C65 C66	120.1	C153 - C154 - II154	120.3
C(4 - C(5 - U(5	120.8 (5)	C154 - C155 - C156	120.1 (5)
C64—C65—H65	119.6	C154—C155—H155	120.0
C66—C65—H65	119.6	C156—C155—H155	120.0
C61—C66—C65	120.1 (3)	C155—C156—C151	121.1 (3)
C61—C66—H66	119.9	C155—C156—H156	119.5
С65—С66—Н66	119.9	C151—C156—H156	119.5
C72—C67—C68	119.4 (3)	C162—C157—C158	119.3 (3)
C72—C67—P3	118.5 (2)	C162—C157—P7	119.1 (2)
C68—C67—P3	122.1 (2)	C158—C157—P7	121.5 (2)
C67—C68—C69	120.6 (3)	C159—C158—C157	120.0 (3)
С67—С68—Н68	119.7	C159—C158—H158	120.0
С69—С68—Н68	119.7	C157—C158—H158	120.0
C70—C69—C68	119.3 (3)	C160—C159—C158	120.2 (3)
С70—С69—Н69	120.4	C160—C159—H159	119.9
С68—С69—Н69	120.4	C158—C159—H159	119.9
С71—С70—С69	120.1 (3)	C161—C160—C159	119.8 (3)
С71—С70—Н70	119.9	C161—C160—H160	120.1
С69—С70—Н70	119.9	C159—C160—H160	120.1
C70—C71—C72	120.7 (3)	C160—C161—C162	120.7 (3)
C70—C71—H71	1197	C160—C161—H161	119.6
C72-C71-H71	119.7	$C_{162} - C_{161} - H_{161}$	119.6
C67 - C72 - C71	119.7	C_{157} $-C_{162}$ $-C_{161}$	119.0 119.9(3)
C67 C72 H72	119.9 (3)	C157 = C162 = C161	119.9 (3)
$C_{0}^{-1} = C_{12}^{-11/2}$	120.1	$C_{157} = C_{162} = H_{162}$	120.0
C74 $C72$ $C78$	120.1	$C_{101} = C_{102} = 11102$	120.0 118.7(2)
C74 - C72 - D4	119.0 (3)	C104 - C103 - C108	118.7(3)
$C_{14} - C_{13} - C_{14}$	110.3(2)	$C104 - C103 - F\delta$	117.9(2)
$C_{10} - C_{10} - C$	122.3(2)	C105 - C103 - P8	125.5(2)
$C_{13} - C_{14} - C_{13}$	120.2 (3)	C103 - C104 - C103	121.2 (3)
C/5—C/4—H/4	119.9	C165—C164—H164	119.4
С/3—С/4—Н74	119.9	C163—C164—H164	119.4
C76—C75—C74	120.1 (3)	C166—C165—C164	119.4 (3)
С76—С75—Н75	120.0	C166—C165—H165	120.3
С74—С75—Н75	120.0	C164—C165—H165	120.3

C75—C76—C77	120.5 (3)	C167—C166—C165	120.1 (3)
С75—С76—Н76	119.7	C167—C166—H166	119.9
С77—С76—Н76	119.7	C165—C166—H166	119.9
C76—C77—C78	119.8 (3)	C166—C167—C168	120.3 (3)
С76—С77—Н77	120.1	C166—C167—H167	119.8
С78—С77—Н77	120.1	C168—C167—H167	119.8
C77—C78—C73	120.4 (3)	C163—C168—C167	120.2 (3)
С77—С78—Н78	119.8	C163—C168—H168	119.9
С73—С78—Н78	119.8	C167—C168—H168	119.9
C80—C79—C84	119.1 (3)	C174—C169—C170	119.3 (3)
C80—C79—P4	123.4 (2)	C174—C169—P8	120.7(3)
C84—C79—P4	117.5 (2)	C170—C169—P8	119.8 (3)
C79 - C80 - C81	1199(3)	$C_{169} - C_{170} - C_{171}$	119.6(3)
C79 - C80 - H80	120.0	$C_{169} - C_{170} - H_{170}$	120.2
C81 - C80 - H80	120.0	$C_{171} - C_{170} - H_{170}$	120.2
C82 - C81 - C80	120.0 120.4(3)	C172 - C171 - C170	120.2 120.0(4)
C82 - C81 - H81	119.8	C_{172} C_{171} H_{171}	120.0 (1)
C80 - C81 - H81	119.8	C170 - C171 - H171	120.0
C_{81} C_{82} C_{83}	119.8 (3)	C173 - C172 - C171	120.0 120.2(4)
$C_{81} C_{82} C_{83}$	120.1	C173 - C172 - C171	110.0
$C_{83} = C_{82} = H_{82}$	120.1	C173 - C172 - H172	119.9
$C_{82} = C_{82} = C$	120.1 120.4(3)	C172 - C172 - C174	119.9
$C_{82} = C_{83} = C_{84}$	110.8	C172 - C173 - C174	120.3(3)
$C_{02} = C_{03} = 1103$	119.0	C172 - C173 - H173	119.7
$C_{84} = C_{83} = H_{83}$	119.0	$C_{1/4} = C_{1/3} = H_{1/3}$	119.7
$C_{83} = C_{84} = C_{79}$	120.5 (5)	C169 - C174 - C173	120.3 (4)
$C_{0} C_{0} C_{0} U_{0} U_{0}$	119.0	C172 - C174 - H174	119.9
C/9 = C84 = H84	119.6	C175 - C174 - H174	119.9
C90 = C85 = C80	110.4(3) 122.1(2)	C176 - C175 - D8	118.9(3)
$C_{90} = C_{85} = P_{4}$	123.1(2)	C170 - C175 - P8	117.0(2)
C80 - C85 - P4	118.5 (2)	C180 - C175 - P8	123.4 (3)
C87 - C86 - C85	120.3 (3)	C1/5 - C1/6 - C1/7	120.9 (3)
C87 - C86 - H86	119.8	C1/5 - C1/6 - H1/6	119.6
C85—C86—H86	119.8	C17/-C176-H176	119.6
	120.4 (3)	C1/8 - C1/7 - C1/6	119.5 (3)
C88—C87—H87	119.8	C178—C177—H177	120.2
C86—C8/—H8/	119.8	C1/6—C1//—H1//	120.2
	120.2 (3)	C179—C178—C177	119.8 (3)
С87—С88—Н88	119.9	C179—C178—H178	120.1
С89—С88—Н88	119.9	C177—C178—H178	120.1
C88—C89—C90	120.1 (3)	C178—C179—C180	121.5 (3)
С88—С89—Н89	119.9	C178—C179—H179	119.3
C90—C89—H89	119.9	С180—С179—Н179	119.3
C89—C90—C85	120.5 (3)	C179—C180—C175	119.4 (3)
С89—С90—Н90	119.7	C179—C180—H180	120.3
С85—С90—Н90	119.7	C175—C180—H180	120.3
C1—N1—N2—C2	-175.1 (3)	C91—N11—N12—C92	175.4 (3)
P2—Ag1—P1—C16	171.02 (11)	P6—Ag3—P5—C106	45.33 (12)

Cl1—Ag1—P1—C16	-42.66 (11)	Cl3—Ag3—P5—C106	-99.22 (12)
S1—Ag1—P1—C16	50.32 (11)	S3—Ag3—P5—C106	166.21 (12)
P2—Ag1—P1—C10	-65.75 (12)	P6—Ag3—P5—C112	-71.79 (10)
Cl1—Ag1—P1—C10	80.56 (12)	Cl3—Ag3—P5—C112	143.66 (10)
S1—Ag1—P1—C10	173.54 (12)	S3—Ag3—P5—C112	49.10 (10)
P2—Ag1—P1—C22	54.75 (11)	P6 - Ag3 - P5 - C100	168.89 (10)
$C_1 - Ag_1 - P_1 - C_{22}$	-158.94(10)	C13 - Ag3 - P5 - C100	24.33 (11)
S1 - Ag1 - P1 - C22	-65.95 (11)	$S_{3} = A_{g_{3}} = P_{5} = C_{100}$	-70.23(11)
P1 - Ag1 - P2 - C34	-72.92(12)	P5 - Ag3 - P6 - C130	-45.98(11)
C11 - Ag1 - P2 - C34	13947(12)	C13 - Ag3 - P6 - C130	100.82(11)
S1 - Ag1 - P2 - C34	40.56(12)	S_{3} Ag3 P_{6} C130	-15884(10)
$P1_{A_{g1}} P2_{C28}$	162.91 (12)	P5 - Ag3 - P6 - C118	-167.00(12)
$C_{11} = A_{g1} = P_{2} = C_{28}$	152.91(12)	$C_{13}^{13} = A_{03}^{10} = P_{6}^{110} = C_{118}^{110}$	-20.20(12)
$S1_Ag1_P2_C28$	-83.61(12)	S_{3} A_{g3} P_{6} C_{118}	80.13(12)
$P1_{Ag1} P2_{C20}$	44 35 (12)	$P5_{4}g3_{7}P6_{124}$	70.80(12)
$\begin{array}{cccccccccccccccccccccccccccccccccccc$	-103.25(12)	$C_{13}^{13} A_{23}^{13} B_{10}^{10} C_{124}^{124}$	-142.30(12)
C_{11} Ag1 B2 C40	103.23(12) 157.84(12)	C13 - Ag3 - 10 - C124	-42.06(12)
S1 - Ag1 - F2 - C40	157.04(12) 162.42(11)	$35 - Ag_3 - F_0 - C_{124}$	-42.00(12)
$P_1 - Ag_1 - S_1 - C_1$	102.42(11) 20.57(11)	$P_{3} = Ag_{3} = S_{3} = C_{9}$	-107.01(11)
P2 - Ag1 - S1 - C1	59.57(11)	P0 - Ag3 - S3 - C91	-42.04(12)
CII - AgI - SI - CI	-77.75(11)	C13 - Ag3 - S3 - C91	$\frac{12.11(11)}{170.2(2)}$
$C_8 N_2 C_1 S_1$	1/7.1(5)	C98 = N13 = C91 = N11	1/9.3 (3)
$C_{0} = N_{0} = C_{1} = S_{1}$	-3.0(4)	C98 - N13 - C91 - S3	-0.4(4)
N2—NI—CI—N3	/.9 (4)	N12—N11—C91—N13	-9.7 (4)
N2—NI—CI—SI	-172.1(2)	N12— $N11$ — $C91$ — $S3$	170.0 (2)
AgI—SI—CI—N3	-130.6 (3)	Ag3—S3—C91—N13	135.6 (2)
Agl—Sl—Cl—Nl	49.3 (3)	Ag3—S3—C91—N11	-44.0 (3)
N1—N2—C2—C4	-178.4 (3)	N11—N12—C92—C94	177.2 (3)
N1—N2—C2—C3	2.9 (5)	N11—N12—C92—C93	-2.0 (5)
C5—N4—C4—C7	4.2 (5)	C95—N14—C94—C97	-3.8 (5)
C5—N4—C4—C2	-174.9 (3)	C95—N14—C94—C92	177.1 (3)
N2—C2—C4—N4	166.8 (3)	N12—C92—C94—N14	-168.3 (3)
C3—C2—C4—N4	-14.5 (5)	C93—C92—C94—N14	11.0 (4)
N2—C2—C4—C7	-12.3 (5)	N12—C92—C94—C97	12.6 (4)
C3—C2—C4—C7	166.4 (3)	C93—C92—C94—C97	-168.2 (3)
C4—N4—C5—C6	-0.7 (6)	C94—N14—C95—C96	0.1 (5)
C7—N5—C6—C5	3.1 (5)	C97—N15—C96—C95	-1.9 (5)
N4—C5—C6—N5	-3.1 (6)	N14—C95—C96—N15	2.9 (6)
C6—N5—C7—C4	0.5 (5)	C96—N15—C97—C94	-1.8 (5)
N4—C4—C7—N5	-4.3 (5)	N14—C94—C97—N15	4.9 (5)
C2—C4—C7—N5	174.8 (3)	C92—C94—C97—N15	-176.1 (3)
C1—N3—C8—C9	-77.7 (4)	C91—N13—C98—C99	79.4 (4)
C16—P1—C10—C15	-32.8 (3)	C106—P5—C100—C105	136.8 (3)
C22—P1—C10—C15	75.4 (3)	C112—P5—C100—C105	-115.4 (3)
Ag1—P1—C10—C15	-160.8 (2)	Ag3—P5—C100—C105	4.2 (3)
C16—P1—C10—C11	149.4 (2)	C106—P5—C100—C101	-40.5 (3)
C22—P1—C10—C11	-102.3 (2)	C112—P5—C100—C101	67.3 (3)
Ag1—P1—C10—C11	21.5 (3)	Ag3—P5—C100—C101	-173.0 (2)
C15-C10-C11-C12	-2.0 (4)	C105—C100—C101—C102	2.5 (4)

P1-C10-C11-C12	175.7 (2)	P5-C100-C101-C102	179.7 (2)
C10-C11-C12-C13	-0.3 (4)	C100-C101-C102-C103	-0.8(4)
C11—C12—C13—C14	2.1 (4)	C101—C102—C103—C104	-1.8(5)
C12—C13—C14—C15	-1.5 (5)	C102—C103—C104—C105	2.6 (5)
C11—C10—C15—C14	2.6 (4)	C101—C100—C105—C104	-1.7(5)
P1-C10-C15-C14	-175.1 (2)	P5-C100-C105-C104	-179.0(3)
C13—C14—C15—C10	-0.9 (4)	C103—C104—C105—C100	-0.8 (6)
C10—P1—C16—C21	-84.9(3)	C112—P5—C106—C107	179.9 (2)
C_{22} P1 $-C_{16}$ C21	167.7 (3)	C100 - P5 - C106 - C107	-70.2(2)
$Ag1_{P1}_{C16}_{C21}$	48.1 (3)	Ag3—P5—C106—C107	61.1 (3)
C10 - P1 - C16 - C17	96.7 (3)	C112—P5—C106—C111	0.0(3)
C_{22} P1 $-C_{16}$ $-C_{17}$	-10.8(3)	C100 - P5 - C106 - C111	109.9(2)
$Ag1_P1_C16_C17$	-1303(2)	$A_{\sigma 3}$ _P5C106_C111	-1188(2)
C_{21} C_{16} C_{17} C_{18}	150.5(2)	$C_{111} - C_{106} - C_{107} - C_{108}$	-0.2(4)
$P_1 = C_16 = C_17 = C_18$	1800(2)	P_{5} C_{106} C_{107} C_{108}	180.0(2)
$C_{16} - C_{17} - C_{18} - C_{19}$	100.0(2)	$C_{106} - C_{107} - C_{108} - C_{109}$	-1.3(4)
$C_{10} = C_{11} = C_{10} = C_{10}$	-1.7(5)	$C_{100} = C_{100} = C_{100} = C_{100}$	20(4)
C18 C19 C20 C21	-0.4(6)	$C_{108} = C_{108} = C_{109} = C_{110}$	-1.3(4)
$C_{10} = C_{10} = C_{20} = C_{21}$	-3.6(5)	$C_{100} = C_{100} = C_{110} = C_{110} = C_{110}$	-0.1(4)
$P_1 = C_{16} = C_{21} = C_{20}$	3.0(3)	$C_{103} = C_{110} = C_{111} = C_{100}$	0.1(4)
11 - 10 - 21 - 220	1/7.9(5)	P5 C106 C111 C110	-170.3(2)
$C_{16} = C_{20} = C_{21} = C_{10}$	-74.5(3)	$C_{106} = C_{100} = C_{111} = C_{110}$	-027(2)
$C_{10} = 1 = C_{22} = C_{27}$	(4.3)	$C_{100} = 15 = C_{112} = C_{117}$	$\frac{92.7(2)}{161.2(2)}$
10-r = -22 - 27	170.3(2)	$100 - r_{3} - C_{112} - C_{117}$	101.5(2)
$Ag_{1} - r_{1} - c_{22} - c_{27}$	47.3(3)	$Ag_{5} - r_{5} - C_{112} - C_{117}$	34.0(2) 87.2(2)
$C_{10} = P_1 = C_{22} = C_{23}$	100.1(3)	$C100 - F_{3} - C112 - C113$	07.2(3)
C10 - P1 - C22 - C23	0.9(3)	C100 - P3 - C112 - C113	-16.6(3)
Ag1 - F1 - C22 - C23	-130.1(2)	Ag_{3} $-r_{3}$ $-c_{112}$ $-c_{113}$	-143.4(2)
$C_2 = C_2 $	2.0(5)	CII/-CII2-CII3-CII4	0.9(4)
P1 - C22 - C23 - C24	1/9.4 (3)	P5—C112—C113—C114	-1/9.0(2)
$C_{22} = C_{23} = C_{24} = C_{25}$	-0.5(5)		0.1(5)
C_{23} C_{24} C_{25} C_{26}	-0.2(5)		-0.9(5)
$C_{24} = C_{25} = C_{26} = C_{27}$	-0.6(5)		0.7(5)
$C_{23} = C_{22} = C_{27} = C_{26}$	-2.8 (5)	C113—C112—C117—C116	-1.1(5)
P1—C22—C27—C26	179.6 (2)	P5—C112—C117—C116	178.8 (3)
C_{25} — C_{26} — C_{27} — C_{22}	2.2 (5)	C115—C116—C117—C112	0.3 (5)
C34—P2—C28—C29	17.2 (3)	C130—P6—C118—C123	101.4 (3)
C40—P2—C28—C29	-91.1 (3)	C124—P6—C118—C123	-4.9 (3)
Ag1—P2—C28—C29	144.6 (2)	Ag3—P6—C118—C123	-132.8 (2)
C34—P2—C28—C33	-162.8 (3)	C130—P6—C118—C119	-78.0 (3)
C40—P2—C28—C33	88.9 (3)	C124—P6—C118—C119	175.7 (2)
Ag1—P2—C28—C33	-35.4 (3)	Ag3—P6—C118—C119	47.9 (3)
C33—C28—C29—C30	-0.9 (5)	C123—C118—C119—C120	-2.3 (5)
P2-C28-C29-C30	179.1 (3)	P6—C118—C119—C120	177.1 (3)
C28—C29—C30—C31	-1.4 (5)	C118—C119—C120—C121	0.5 (5)
C29—C30—C31—C32	1.9 (5)	C119—C120—C121—C122	1.5 (5)
C30—C31—C32—C33	-0.2 (6)	C120—C121—C122—C123	-1.8(5)
C31—C32—C33—C28	-2.1 (5)	C119—C118—C123—C122	2.0 (5)
C29—C28—C33—C32	2.6 (5)	P6-C118-C123-C122	-177.3 (3)

P2-C28-C33-C32	-177.4 (3)	C121—C122—C123—C118	0.0 (5)
C28—P2—C34—C35	112.0 (3)	C130—P6—C124—C125	-20.6 (3)
C40—P2—C34—C35	-140.5 (3)	C118—P6—C124—C125	86.7 (3)
Ag1—P2—C34—C35	-17.6 (3)	Ag3—P6—C124—C125	-143.5 (2)
C28—P2—C34—C39	-69.6 (3)	C130—P6—C124—C129	160.0 (2)
C40—P2—C34—C39	37.9 (3)	C118—P6—C124—C129	-92.7 (2)
Ag1—P2—C34—C39	160.8 (2)	Ag3—P6—C124—C129	37.1 (3)
C39—C34—C35—C36	-1.3 (5)	C129—C124—C125—C126	0.2 (4)
P2-C34-C35-C36	177.2 (3)	P6-C124-C125-C126	-179.2 (2)
C34—C35—C36—C37	0.9 (6)	C124—C125—C126—C127	0.8 (3)
C35—C36—C37—C38	-0.4(5)	C125—C126—C127—C128	-1.1(2)
C36—C37—C38—C39	0.4 (5)	C126—C127—C128—C129	0.4 (4)
C37—C38—C39—C34	-0.8(5)	C125—C124—C129—C128	-1.0(4)
C35—C34—C39—C38	1.2 (5)	P6-C124-C129-C128	178.5 (2)
P2-C34-C39-C38	-177.2(2)	C127—C128—C129—C124	0.6 (5)
C_{34} P2 C40 C41	-150.4(3)	C118—P6—C130—C131	2.2(3)
$C_{28} = P_{2} = C_{40} = C_{41}$	-39.8(3)	C124—P6—C130—C131	109.4 (3)
$Ag1_{P2}_{C40}_{C41}$	86.9 (3)	Ag3—P6—C130—C131	-126.6(2)
C_{34} P2 C40 C45	35.7 (3)	C118—P6—C130—C135	-178.9(2)
C_{28} P2 C_{40} C45	1463(3)	C_{124} P6 C_{130} C_{135}	-71.8(2)
$Ag1_{P2}_{C40}_{C45}$	-87.0(3)	Ag3—P6—C130—C135	52.2(2)
C45-C40-C41-C42	0.6 (5)	C_{135} $-C_{130}$ $-C_{131}$ $-C_{132}$	0.7 (5)
P2-C40-C41-C42	-173.4(3)	P6-C130-C131-C132	179.6 (3)
C40-C41-C42-C43	-1.6(6)	C_{130} $-C_{131}$ $-C_{132}$ $-C_{133}$	0.1 (5)
C41 - C42 - C43 - C44	0.9 (6)	C_{131} $-C_{132}$ $-C_{133}$ $-C_{134}$	-0.8(5)
C42-C43-C44-C45	0.9 (6)	C_{132} C_{133} C_{134} C_{135}	0.6(5)
C41 - C40 - C45 - C44	1.2(5)	C_{133} C_{134} C_{135} C_{130}	0.3(5)
P2-C40-C45-C44	175 2 (3)	$C_{131} - C_{130} - C_{135} - C_{134}$	-0.9(4)
C43 - C44 - C45 - C40	-2.0(6)	P6-C130-C135-C134	-179.8(2)
C46—N6—N7—C47	-176.4(3)	$C_{136} N_{16} N_{17} C_{137}$	171.8 (3)
P4 - Ag2 - P3 - C61	-45.67(12)	P8 - Ag4 - P7 - C145	-170.74(11)
Cl2—Ag2—P3—C61	99.11 (12)	C_{14} Ag4 P_{7} C_{145}	44.22 (11)
S_{2} A_{g2} P_{3} C_{61}	-167.51(12)	S4—Ag4—P7—C145	-49.88(11)
P4 - Ag2 - P3 - C55	71.71 (10)	P8 - Ag4 - P7 - C157	65.57 (13)
C12—Ag2—P3—C55	-143.51(10)	C14—Ag4—P7—C157	-79.47(13)
S2—Ag2—P3—C55	-50.13(10)	S4—Ag4—P7—C157	-173.58(12)
P4 - Ag2 - P3 - C67	-169.38(11)	P8 - Ag4 - P7 - C151	-55.65(11)
C12 - Ag2 - P3 - C67	$-24\ 60\ (12)$	$C_{14} - A_{\sigma 4} - P_{7} - C_{151}$	159 31 (10)
$S_{2}^{2} = A_{3}^{2} = P_{3}^{2} = C_{6}^{2}$	68 79 (11)	S4 - Ag4 - P7 - C151	65 21 (11)
$P_3 = A_{\sigma}^2 = P_4 = C_{85}^{-1}$	-71.76(12)	P7 - Ag4 - P8 - C163	-159 19 (12)
C12 - Ag2 - P4 - C85	141 31 (12)	$C_{14} - A_{\sigma 4} - P_{8} - C_{163}$	-1355(12)
$S_{2}^{2} - A_{9}^{2} - P_{4}^{2} - C_{85}^{2}$	40.90(12)	S4 - Ag4 - P8 - C163	85.09(12)
$P_3 = A_{\sigma}^2 = P_4 = C_{73}^2$	166 14 (12)	P7 - Ag4 - P8 - C169	-40.81(13)
C12 - Ag2 - P4 - C73	19.21(12)	$C_{14} - A_{\sigma 4} - P_{8} - C_{169}$	104 82 (13)
$S_{2} = A_{g2} = P_{4} = C_{73}$	-81.19(12)	S4—Ag4—P8—C169	-156.53(13)
$P_3 = A_{g_2} = P_4 = C_{79}$	44 65 (11)	P7—Ag4—P8—C175	76 10 (12)
C12—Ag2—P4—C79	-102.28(10)	C_{14} Ag4 P8 C175	-138 26 (12)
$S_{2}^{2} = A_{\sigma}^{2} = P_{4}^{2} = C_{7}^{2}$	157 32 (10)	$S4_Ag4_B8_C175$	-39.62(12)
52 Mg2 17 C/3	157.52 (10)	57 ngt 10 C1/5	57.02 (12)

P3—Ag2—S2—C46	167.73 (11)	P7—Ag4—S4—C136	-163.09 (11)
P4—Ag2—S2—C46	41.97 (11)	P8—Ag4—S4—C136	-40.74 (11)
Cl2—Ag2—S2—C46	-72.89 (11)	Cl4—Ag4—S4—C136	77.76 (11)
C53—N8—C46—N6	179.8 (3)	C143—N18—C136—N16	-172.0(3)
C53—N8—C46—S2	-0.9 (4)	C143—N18—C136—S4	7.9 (5)
N7—N6—C46—N8	10.5 (4)	N17—N16—C136—N18	-7.2 (4)
N7—N6—C46—S2	-168.8(2)	N17—N16—C136—S4	172.9 (2)
Ag2-S2-C46-N8	-136.1 (2)	Ag4—S4—C136—N18	133.4 (3)
Ag2—S2—C46—N6	43.2 (3)	Ag4—S4—C136—N16	-46.7(3)
N6—N7—C47—C49	-177.6 (3)	N16—N17—C137—C139	176.8 (3)
N6—N7—C47—C48	2.1 (5)	N16—N17—C137—C138	-2.3 (5)
C50—N9—C49—C52	4.3 (5)	C140—N19—C139—C142	-4.9(5)
C50—N9—C49—C47	-173.5 (3)	C140—N19—C139—C137	174.4 (3)
N7—C47—C49—N9	166.7 (3)	N17—C137—C139—N19	-163.8(3)
C48—C47—C49—N9	-13.0 (4)	C138—C137—C139—N19	15.4 (4)
N7—C47—C49—C52	-11.2(4)	N17—C137—C139—C142	15.5 (4)
C48—C47—C49—C52	169.2 (3)	C138—C137—C139—C142	-165.3(3)
C49—N9—C50—C51	0.4 (5)	C139—N19—C140—C141	0.4 (6)
C52—N10—C51—C50	3.8 (5)	C142—N20—C141—C140	-2.5(6)
N9-C50-C51-N10	-4.7(6)	N19—C140—C141—N20	3.5(7)
C51—N10—C52—C49	1.0 (5)	C141 - N20 - C142 - C139	-2.1(5)
N9-C49-C52-N10	-5.3(5)	N19—C139—C142—N20	6.0 (5)
C47-C49-C52-N10	172.5 (3)	C_{137} C_{139} C_{142} N_{20}	-1733(3)
C46 - N8 - C53 - C54	-801(4)	C136—N18— $C143$ — $C144$	86 4 (4)
C61 - P3 - C55 - C56	93 1 (3)	C157 - P7 - C145 - C146	-90.1(3)
C67 - P3 - C55 - C56	-1605(2)	C151 - P7 - C145 - C146	171(3)
Ag2 - P3 - C55 - C56	-351(3)	Ag4 - P7 - C145 - C146	1354(2)
C_{61} P_{3} C_{55} C_{60}	-87.9(3)	C_{157} P7 C_{145} C150	90.1 (3)
C67 - P3 - C55 - C60	18 5 (3)	C_{151} P7- C_{145} C150	-162.6(3)
Ag2 - P3 - C55 - C60	143.9(2)	Ag4 = P7 = C145 = C150	-444(3)
C60 - C55 - C56 - C57	21(5)	C_{150} C_{145} C_{146} C_{147}	-24(5)
P3-C55-C56-C57	-1788(3)	P7—C145—C146—C147	177.8(2)
$C_{55} - C_{56} - C_{57} - C_{58}$	-0.8(5)	$C_{145} = C_{146} = C_{147} = C_{148}$	-0.1(5)
$C_{56} - C_{57} - C_{58} - C_{59}$	-0.8(5)	C146 $C147$ $C148$ $C149$	15(5)
C_{57} C_{58} C_{59} C_{60}	11(5)	C147 - C148 - C149 - C150	-0.4(6)
C_{58} C_{59} C_{60} C_{55}	0.2(5)	$C_{148} = C_{149} = C_{150} = C_{145}$	-22(6)
$C_{56} - C_{55} - C_{60} - C_{59}$	-1.8(4)	C146 - C145 - C150 - C149	36(5)
P3-C55-C60-C59	179 3 (2)	P7-C145-C150-C149	-1767(3)
C_{55} P_{3} C_{61} C_{66}	-179.7(2)	$C_{145} = P_{7} = C_{151} = C_{152}$	78 6 (3)
C67 - P3 - C61 - C66	703(2)	C157 - P7 - C151 - C152	-1743(2)
A_{g2} P3 C61 C66	-60.4(2)	A g 4 P 7 C 151 C 152	-414(2)
C_{55} P3 C_{61} C_{62}	-1.2(3)	C_{145} P7 C_{151} C_{152}	-1035(3)
C67 - P3 - C61 - C62	-1113(2)	$C_{145} = P_{7} = C_{151} = C_{150}$	37(3)
A_{g2} P3 C61 C62	111.3(2) 118.0(2)	$\Delta q A = P7 = C151 = C156$	3.7(3)
$\Gamma_{66} = \Gamma_{61} = \Gamma_{61} = \Gamma_{62} = \Gamma_{63}$	-1.6(4)	C_{156} C_{151} C_{152} C_{152}	150.0(2) 15(4)
$P_3 = C_{61} = C_{62} = C_{63}$	-1800(2)	P7 C151 C152 C153	1.3(4) 1706(2)
$C_{61} C_{62} C_{62} C_{64}$	100.0(2)	17 - 0151 - 0152 - 0155	-0.7(5)
$C_{01} = C_{02} = C_{03} = C_{04}$	0.3(4)	$C_{131} - C_{132} - C_{133} - C_{134}$	-0.3(5)
02 - 03 - 04 - 03	1.0 (4)	U132 - U133 - U134 - U133	-0.5 (3)

C63—C64—C65—C66	-1.4 (4)	C153—C154—C155—C156	0.3 (5)
C62—C61—C66—C65	1.2 (4)	C154—C155—C156—C151	0.6 (5)
P3-C61-C66-C65	179.7 (2)	C152—C151—C156—C155	-1.5(5)
C64—C65—C66—C61	0.3 (4)	P7—C151—C156—C155	-179.5 (2)
C61—P3—C67—C72	-136.3 (3)	C145—P7—C157—C162	-149.3 (2)
C55—P3—C67—C72	116.0 (3)	C151—P7—C157—C162	102.5 (2)
Ag2—P3—C67—C72	-3.3 (3)	Ag4—P7—C157—C162	-21.6(3)
C61—P3—C67—C68	40.0 (3)	C145—P7—C157—C158	33.4 (3)
C55 - P3 - C67 - C68	-67.6 (3)	C151—P7—C157—C158	-74.8(3)
Ag2—P3—C67—C68	173.1 (2)	Ag4—P7—C157—C158	161.1 (2)
C72-C67-C68-C69	-36(4)	C_{162} C_{157} C_{158} C_{159}	-32(4)
P_{3} C_{67} C_{68} C_{69}	-1799(2)	P7-C157-C158-C159	1742(2)
C67 - C68 - C69 - C70	16(4)	C_{157} C_{158} C_{159} C_{160}	171.2(2) 13(4)
C68 - C69 - C70 - C71	0.8(5)	C_{158} C_{159} C_{160} C_{161}	1.3(+) 1 2 (4)
C69 $C70$ $C71$ $C72$	-12(5)	$C_{150} = C_{150} = C_{160} = C_{161} = C_{161}$	-1.8(5)
C68 C67 C72 C71	1.2(3)	$C_{159} = C_{100} = C_{101} = C_{102}$	26(4)
$C_{00} = C_{01} = C_{12} = C_{11}$	5.1(5)	P7 C157 C162 C161	2.0(4) -174.8(2)
$r_{3} = c_{0} = c_{1} = c_{1}$	1/9.3(3) -0.7(5)	$\Gamma = C157 = C102 = C101$	-1/4.0(2) -0.2(4)
$C^{95} P_{4} C^{72} C^{74}$	-0.7(3)	$C_{100} = C_{101} = C_{102} = C_{157}$	-0.2(4)
C_{83} P4 C_{73} C_{74}	-1/4.1(3)	C109 - P8 - C103 - C104	-88.0(3)
C/9 - P4 - C/3 - C/4	/9.4 (3)	C1/3 - P8 - C103 - C104	103.0(3)
Ag2 - P4 - C73 - C74	-40.8(3)	Ag4 - P8 - C103 - C104	54.8 (5) 00.5 (2)
C_{85} P4 C_{73} C_{78}	6.1(3)	C169 - P8 - C163 - C168	90.5 (3)
C/9 - P4 - C/3 - C/8	-100.4(3)	C1/5 - P8 - C163 - C168	-1/.9(3)
Ag2—P4—C/3—C/8	133.4 (2)	Ag4—P8—C163—C168	-146.1 (2)
C/8—C/3—C/4—C/5	3.1 (5)	C168—C163—C164—C165	-1.8(5)
P4—C73—C74—C75	-176.7 (3)	P8—C163—C164—C165	177.4 (3)
C73—C74—C75—C76	-2.5 (5)	C163—C164—C165—C166	1.7 (5)
C74—C75—C76—C77	0.1 (5)	C164—C165—C166—C167	-0.5(5)
C75—C76—C77—C78	1.7 (5)	C165—C166—C167—C168	-0.5(5)
C76—C77—C78—C73	-1.1 (5)	C164—C163—C168—C167	0.8 (5)
C74—C73—C78—C77	-1.3 (5)	P8—C163—C168—C167	-178.3 (3)
P4—C73—C78—C77	178.5 (2)	C166—C167—C168—C163	0.3 (5)
C85—P4—C79—C80	-109.9 (3)	C163—P8—C169—C174	36.9 (3)
C73—P4—C79—C80	-2.4 (3)	C175—P8—C169—C174	147.6 (3)
Ag2—P4—C79—C80	127.0 (2)	Ag4—P8—C169—C174	-89.6 (3)
C85—P4—C79—C84	71.8 (2)	C163—P8—C169—C170	-148.7 (3)
C73—P4—C79—C84	179.3 (2)	C175—P8—C169—C170	-38.0 (3)
Ag2—P4—C79—C84	-51.3 (2)	Ag4—P8—C169—C170	84.8 (3)
C84—C79—C80—C81	-0.2 (4)	C174—C169—C170—C171	-1.6 (5)
P4-C79-C80-C81	-178.5 (2)	P8-C169-C170-C171	-176.2 (3)
C79—C80—C81—C82	-1.2 (5)	C169—C170—C171—C172	2.4 (6)
C80—C81—C82—C83	1.6 (5)	C170—C171—C172—C173	-2.2 (7)
C81—C82—C83—C84	-0.5 (5)	C171—C172—C173—C174	1.1 (7)
C82—C83—C84—C79	-1.0 (5)	C170—C169—C174—C173	0.6 (5)
C80—C79—C84—C83	1.3 (5)	P8—C169—C174—C173	175.1 (3)
P4C79C84C83	179.7 (3)	C172—C173—C174—C169	-0.4 (6)
C73—P4—C85—C90	-84.4 (3)	C163—P8—C175—C176	-109.5 (3)
C79—P4—C85—C90	23.2 (3)	C169—P8—C175—C176	142.6 (3)

A ~2 D4 C95 C00	145.9 (2)	A = 4 D9 C175 C176	20.7(2)
Ag2—P4—C85—C90	145.8 (2)	Ag4—P8—C1/5—C1/6	20.7 (3)
C73—P4—C85—C86	95.2 (2)	C163—P8—C175—C180	73.0 (3)
C79—P4—C85—C86	-157.3 (2)	C169—P8—C175—C180	-34.9 (3)
Ag2—P4—C85—C86	-34.6 (3)	Ag4—P8—C175—C180	-156.8 (2)
C90—C85—C86—C87	-0.4 (4)	C180—C175—C176—C177	0.0 (5)
P4—C85—C86—C87	-179.9 (2)	P8—C175—C176—C177	-177.6 (3)
C85—C86—C87—C88	0.6 (5)	C175—C176—C177—C178	1.2 (5)
C86—C87—C88—C89	-0.6 (5)	C176—C177—C178—C179	-1.9 (5)
C87—C88—C89—C90	0.4 (5)	C177—C178—C179—C180	1.4 (5)
C88—C89—C90—C85	-0.2 (5)	C178—C179—C180—C175	-0.1 (5)
C86—C85—C90—C89	0.2 (4)	C176—C175—C180—C179	-0.6 (5)
P4—C85—C90—C89	179.8 (2)	P8—C175—C180—C179	176.9 (2)

Hydrogen-bond geometry (Å, °)

D—H···A	<i>D</i> —Н	H···A	$D \cdots A$	D—H···A
N1—H1 <i>N</i> …Cl1	0.88 (2)	2.54 (2)	3.381 (3)	163 (3)
N3—H3 <i>N</i> ···Cl3	0.87 (2)	2.68 (3)	3.406 (3)	142 (3)
N6—H6 <i>N</i> ···Cl2	0.88 (2)	2.53 (2)	3.370 (3)	162 (3)
N8—H8 <i>N</i> ···Cl4	0.86 (2)	2.68 (3)	3.411 (3)	143 (3)
N11—H11 <i>N</i> ···Cl3	0.88 (2)	2.52 (2)	3.363 (3)	161 (3)
N13—H13 <i>N</i> ···Cl2	0.88 (2)	2.65 (2)	3.423 (3)	147 (3)
N16—H16 <i>N</i> ···Cl4	0.86 (2)	2.50 (2)	3.338 (3)	164 (3)
N18—H18N····Cl1 ⁱ	0.86 (2)	2.70 (3)	3.425 (3)	143 (3)

Symmetry code: (i) x-1, y, z-1.