

 $\nu = 90.383 \ (1)^{\circ}$ 

Z = 4

V = 4692.38 (19) Å<sup>3</sup>

 $0.35 \times 0.23 \times 0.19 \text{ mm}$ 

161249 measured reflections 40931 independent reflections

35153 reflections with  $I > 2\sigma(I)$ 

Mo  $K\alpha$  radiation

 $\mu = 2.45 \text{ mm}^{-1}$ 

T = 100 K

 $R_{\rm int} = 0.026$ 

CH

CH

CH.

Acta Crystallographica Section E Structure Reports Online

ISSN 1600-5368

# [ $\mu$ -Bis(diphenylarsanyl)methane-1:2 $\kappa^2$ As:As']nonacarbonyl-1 $\kappa^3$ C,2 $\kappa^3$ C,3 $\kappa^3$ C-(triisopropyl phosphite-3 $\kappa$ P)-triangulo-triruthenium(0)

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Received 19 March 2011; accepted 31 March 2011

Key indicators: single-crystal X-ray study; T = 100 K; mean  $\sigma$ (C–C) = 0.002 Å; R factor = 0.024; wR factor = 0.055; data-to-parameter ratio = 37.4.

The asymmetric unit of the title *triangulo*-triruthenium compound, Ru<sub>3</sub>(CO)<sub>9</sub>( $\mu$ -Ph<sub>2</sub>AsCH<sub>2</sub>AsPh<sub>2</sub>)(P[OCH(CH<sub>3</sub>)<sub>2</sub>]<sub>3</sub>) or [Ru<sub>3</sub>(C<sub>25</sub>H<sub>22</sub>As<sub>2</sub>)(C<sub>9</sub>H<sub>21</sub>O<sub>3</sub>P)(CO)<sub>9</sub>], contains two molecules of the *triangulo*-triruthenium complex. The bis(diphenylarsanyl)methane ligand bridges an Ru-Ru bond and the monodentate phosphite ligands to the third Ru atom. Both the arsine and phosphite ligands are equatorial with respect to the Ru<sub>3</sub> triangle. Additionally, each Ru atom carries one equatorial and two axial terminal carbonyl ligands. The dihedral angles between the pairs of benzene rings bound to individual As atoms are 85.67 (8) and 75.91 (8) in the first independent molecule and 74.64 (8) and 70.76 (9) in the second. In the crystal, molecules are linked into a three-dimensional framework by intermolecular C-H···O hydrogen bonds.

#### **Related literature**

For related structures, see: Bruce *et al.* (1983, 1988); Churchill *et al.* (1977); Shawkataly *et al.* (1998). For the synthesis, see: Shawkataly *et al.* (2011); Bruce *et al.* (1987). For the stability of the temperature controller used in the data collection, see: Cosier & Glazer (1986).

#### **Experimental**

#### Crystal data

 $[Ru_{3}(C_{25}H_{22}As_{2})(C_{9}H_{21}O_{3}P)(CO)_{9}]$   $M_{r} = 1235.79$ Triclinic,  $P\overline{1}$  a = 12.3481 (3) Å b = 18.1697 (4) Å c = 21.5295 (5) Å  $\alpha = 93.494$  (1)°  $\beta = 103.230$  (1)°

#### Data collection

Bruker APEXII DUO CCD areadetector diffractometer Absorption correction: multi-scan (SADABS; Bruker, 2009)  $T_{min} = 0.483, T_{max} = 0.659$ 

#### Refinement

 $R[F^2 > 2\sigma(F^2)] = 0.024$ 1093 parameters $wR(F^2) = 0.055$ H-atom parameters constrainedS = 1.01 $\Delta \rho_{max} = 1.57$  e Å $^{-3}$ 40931 reflections $\Delta \rho_{min} = -1.26$  e Å $^{-3}$ 

#### Table 1

Hydrogen-bond geometry (Å, °).

$D - H \cdots A$	$D-\mathrm{H}$	$H \cdot \cdot \cdot A$	$D \cdots A$	$D - \mathbf{H} \cdots A$
$C10B - H10B \cdots O6A^{i}$ $C17B - H17B \cdots O2B^{ii}$ $C33A - H33B \cdots O2A^{iii}$	0.93 0.93 0.96	2.59 2.57 2.56	3.293 (2) 3.477 (2) 3.492 (2)	132 165 164
Symmetry codes: (i) -x, -y + 1, -z + 1.	-x, -y + 1	, − <i>z</i> ; (ii)	-x + 1, -y + 2	1, -z; (iii)

Data collection: *APEX2* (Bruker, 2009); cell refinement: *SAINT* (Bruker, 2009); data reduction: *SAINT*; program(s) used to solve structure: *SHELXTL* (Sheldrick, 2008); program(s) used to refine structure: *SHELXTL*; molecular graphics: *SHELXTL*; software used to prepare material for publication: *SHELXTL* and *PLATON* (Spek, 2009).

We gratefully acknowledge funding from the Malaysian Government and Universiti Sains Malaysia (USM) under the University Research Grant 1001/PJJAUH/811115. MGA thanks USM for a dost-doctoral fellowship and HKF thanks the Malaysian Government and Universiti Sains Malaysia for the Research University Grant No. 1001/PFIZIK/811160.

<sup>‡</sup> Thomson Reuters ResearcherID: B-6034-2009.

<sup>§</sup> Thomson Reuters Researcher ID: G-4202-2010.

<sup>¶</sup> Thomson Reuters ResearcherID: A-5523-2009. ‡‡ Thomson Reuters ResearcherID: A-3561-2009.

Supplementary data and figures for this paper are available from the IUCr electronic archives (Reference: SJ5121).

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# supporting information

Acta Cryst. (2011). E67, m545-m546 [doi:10.1107/S1600536811012050]

# [ $\mu$ -Bis(diphenylarsanyl)methane-1:2 $\kappa^2$ As:As']nonacarbonyl-1 $\kappa^3$ C,2 $\kappa^3$ C,3 $\kappa^3$ C-(triisopropyl phosphite-3 $\kappa$ P)-triangulo-triruthenium(0)

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#### S1. Comment

Syntheses and crystal structures of substituted *triangulo*-triruthenium clusters have been of interest to researchers due to structural variations and catalytic activities. A large number of substituted derivatives,  $Ru_3(CO)_{12-n}L_n$  (L = group 15 ligand) have been reported (Bruce *et al.*, 1988). As part of our ongoing studies (Shawkataly *et al.*, 2011), we report here a mixed-ligand metal-carbonyl cluster complex,  $Ru_3(CO)_9(\mu$ -Ph<sub>2</sub>AsCH<sub>2</sub>AsPh<sub>2</sub>)(P[OCH(CH<sub>3</sub>)<sub>2</sub>]<sub>3</sub>).

The asymmetric unit consists of two crystallographically independent molecules (Fig. 1) of the *triangulo*-triruthenium complex, A and B. Fig. 2 shows the atom numbering scheme for molecule A. Atoms of molecule B were labelled similarly. The bond lengths and angles of the title compound are comparable to those found in its related structure (Shawkataly *et al.*, 1998). The bis(diphenylarsino)methane ligand bridges the Ru1 – Ru2 bond and the monodentate phosphite ligand bonds to the Ru3 atom. Both the phosphite and arsine ligands are equatorial with respect to the Ru<sub>3</sub> triangle. Additionally, each Ru atom carries one equatorial and two axial terminal carbonyl ligands. The dihedral angles between the two benzene rings (C1-C6/C7-C12 & C14-C19/C20-C25) are 85.67 (8), 75.91 (8) and 74.64 (8), 70.76 (9) for the two diphenylarsanyl groups of molecule A and B, respectively.

In the title compound, one of the Ru - Ru bonds is noticeably longer compared to the other two Ru - Ru bonds. The unevenness in the lengths of Ru - Ru bonds in comparison with those of the  $Ru_3(CO)_{12}$  structure (Churchill *et al.*, 1977), can be attributed to the steric effect induced by the bulky substituent. In the crystal, molecules are linked into a three-dimensional framework by intermolecular C—H…O weak intearctions. (Fig. 3).

#### **S2. Experimental**

All manipulations were performed under a dry, oxygen-free dinitrogen atmosphere using standard Schlenk techniques. All solvents were dried over sodium and distilled from sodium benzophenone ketyl under nitrogen (Bruce *et al.*, 1987). Isopropyl phosphite was used as received and bis(diphenylarsino)methane (Bruce *et al.*, 1983) was prepared by the reported procedure.

Equimolar quantities of  $Ru_3(CO)_{10}(\mu$ -Ph<sub>2</sub>AsCH<sub>2</sub>AsPh) and isopropyl phosphite were refluxed in THF under nitrogen for 20 minutes. The reaction mixture turned intense red. The solvent was removed under vacuum. The reaction mixture was separated by TLC (dichloromethane:hexane, 1:3). Two bands appeared. The major band (red)  $R_f$ = 0.61 was separated and characterized. The compound was crystallised from CH<sub>2</sub>Cl<sub>2</sub>- CH<sub>3</sub>OH, 1:3 yield = 55.2 %, m.p. 189°C. IR(cyclohexane): v (CO) 2127s, 2059 vs, 2042 s, 2017 s, 1988 s cm<sup>-1</sup>.

#### S3. Refinement

All hydrogen atoms were positioned geometrically and refined using a riding model with C—H = 0.93-0.98 Å and  $U_{iso}(H) = 1.2U_{eq}(C)$  and  $1.5U_{eq}$  (C of methyl group).



### Figure 1

The asymmetric unit, showing 50% probability displacement ellipsoids.



#### Figure 2

The structure of molecule A, showing 50% probability displacement ellipsoids and the atom-numbering scheme. Atoms in molecule B are similarly labelled.



#### Figure 3

The crystal packing of (I) viewed along the c axis. Dashed lines indicate hydrogen bonds. H atoms not involved in the hydrogen bond interactions have been omitted for clarity.

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#### Crystal data

2	
$[Ru_{3}(C_{25}H_{22}As_{2})(C_{9}H_{21}O_{3}P)(CO)_{9}]$ $M_{r} = 1235.79$ Triclinic, $P\overline{1}$ Hall symbol: -P 1 a = 12.3481 (3) Å b = 18.1697 (4) Å c = 21.5295 (5) Å a = 93.494 (1)° $\beta = 103.230$ (1)° $\gamma = 90.383$ (1)° V = 4692.38 (19) Å <sup>3</sup>	Z = 4 F(000) = 2440 $D_x = 1.749 \text{ Mg m}^{-3}$ Mo Ka radiation, $\lambda = 0.71073 \text{ Å}$ Cell parameters from 9785 reflections $\theta = 2.8-37.6^{\circ}$ $\mu = 2.45 \text{ mm}^{-1}$ T = 100 K Plate, red $0.35 \times 0.23 \times 0.19 \text{ mm}$
v = 4022.30 (12) A	
Data collection	
Bruker APEXII DUO CCD area-detector diffractometer Radiation source: fine-focus sealed tube Graphite monochromator $\varphi$ and $\omega$ scans Absorption correction: multi-scan ( <i>SADABS</i> ; Bruker, 2009) $T_{min} = 0.483, T_{max} = 0.659$	161249 measured reflections 40931 independent reflections 35153 reflections with $I > 2\sigma(I)$ $R_{int} = 0.026$ $\theta_{max} = 35.0^\circ, \theta_{min} = 1.8^\circ$ $h = -19 \rightarrow 19$ $k = -29 \rightarrow 29$ $l = -34 \rightarrow 34$
Refinement	
Refinement on $F^2$ Least-squares matrix: full $R[F^2 > 2\sigma(F^2)] = 0.024$ $wR(F^2) = 0.055$	S = 1.01 40931 reflections 1093 parameters 0 restraints

H-atom parameters constrained
$w = 1/[\sigma^2(F_o^2) + (0.0207P)^2 + 3.4447P]$
where $P = (F_o^2 + 2F_c^2)/3$
$(\Delta/\sigma)_{\rm max} = 0.005$
$\Delta \rho_{\rm max} = 1.57 \text{ e } \text{\AA}^{-3}$
$\Delta \rho_{\rm min} = -1.26 \text{ e } \text{\AA}^{-3}$

#### Special details

**Experimental**. The crystal was placed in the cold stream of an Oxford Cyrosystems Cobra open-flow nitrogen cryostat (Cosier & Glazer, 1986) operating at 100.0 (1) K.

**Geometry**. All esds (except the esd in the dihedral angle between two l.s. planes) are estimated using the full covariance matrix. The cell esds are taken into account individually in the estimation of esds in distances, angles and torsion angles; correlations between esds in cell parameters are only used when they are defined by crystal symmetry. An approximate (isotropic) treatment of cell esds is used for estimating esds 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  $F^2 > 2$ 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.

	x	У	Ζ	$U_{ m iso}$ */ $U_{ m eq}$	
Ru1A	0.185622 (9)	0.312340 (6)	0.427544 (5)	0.01368 (2)	
Ru2A	0.340063 (9)	0.193747 (6)	0.432267 (5)	0.01354 (2)	
Ru3A	0.197922 (9)	0.197674 (6)	0.517361 (5)	0.01439 (2)	
As1A	0.236634 (11)	0.355824 (7)	0.331871 (6)	0.01348 (2)	
As2A	0.428726 (11)	0.234754 (8)	0.351165 (6)	0.01387 (2)	
P1A	0.05349 (3)	0.20550 (2)	0.567872 (17)	0.01507 (6)	
O1A	-0.03079 (10)	0.22772 (7)	0.36225 (6)	0.0276 (2)	
O2A	0.03402 (11)	0.43647 (7)	0.44928 (6)	0.0281 (3)	
O3A	0.37695 (11)	0.39986 (7)	0.51876 (6)	0.0268 (2)	
O4A	0.51675 (10)	0.26282 (7)	0.54581 (6)	0.0247 (2)	
O5A	0.43156 (12)	0.03960 (7)	0.44851 (7)	0.0338 (3)	
O6A	0.16098 (10)	0.13352 (7)	0.31667 (6)	0.0269 (2)	
O7A	0.04765 (11)	0.08535 (7)	0.42250 (6)	0.0293 (2)	
O8A	0.32119 (11)	0.06903 (7)	0.58178 (6)	0.0302 (3)	
09A	0.33184 (11)	0.30623 (8)	0.62211 (6)	0.0303 (3)	
O10A	0.04745 (9)	0.14607 (6)	0.61850 (5)	0.01956 (19)	
011A	-0.06140 (9)	0.18840 (6)	0.51830 (5)	0.01920 (19)	
O12A	0.03848 (9)	0.27785 (6)	0.61109 (5)	0.01966 (19)	
C1A	0.12125 (11)	0.38267 (7)	0.25888 (7)	0.0154 (2)	
C2A	0.02752 (13)	0.41811 (9)	0.27021 (8)	0.0230 (3)	
H2AA	0.0163	0.4234	0.3115	0.028*	
C3A	-0.04975 (14)	0.44579 (10)	0.21959 (9)	0.0270 (3)	
H3AA	-0.1118	0.4701	0.2274	0.032*	
C4A	-0.03447 (13)	0.43732 (9)	0.15792 (8)	0.0229 (3)	
H4AA	-0.0861	0.4560	0.1244	0.027*	
C5A	0.05729 (14)	0.40114 (10)	0.14619 (8)	0.0258 (3)	
H5AA	0.0667	0.3945	0.1046	0.031*	
C6A	0.13607 (13)	0.37452 (9)	0.19675 (7)	0.0234 (3)	

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

H6AA	0.1988	0.3512	0.1888	0.028*
C7A	0.32135 (11)	0.44843 (8)	0.34206 (7)	0.0158 (2)
C8A	0.38946 (14)	0.46559 (9)	0.30141 (8)	0.0221 (3)
H8AA	0.4005	0.4307	0.2701	0.027*
C9A	0.44112 (14)	0.53480 (9)	0.30737 (8)	0.0239 (3)
H9AA	0.4866	0.5458	0.2801	0.029*
C10A	0.42536 (13)	0.58731 (9)	0.35355 (8)	0.0222 (3)
H10A	0.4602	0.6334	0.3574	0.027*
C11A	0.35700 (14)	0.57061 (9)	0.39414 (8)	0.0231 (3)
H11A	0.3459	0.6058	0.4252	0.028*
C12A	0.30522 (13)	0.50151 (8)	0.38852 (7)	0.0195 (2)
H12A	0.2597	0.4907	0.4158	0.023*
C13A	0.32281(12)	0.28819(8)	0.28830 (6)	0.0162(2)
H13A	0.2728	0.2533	0.2594	0.0102 (2)
H13R	0.3628	0.3159	0.2632	0.019*
C14A	0.3020 0.47177(12)	0.5155 0.15555(7)	0.2032 0.29804 (7)	0.019
	0.47177(12) 0.54633(13)	0.10333(9)	0.29804(7) 0.32842(7)	0.0103(2) 0.0217(3)
H15A	0.5753	0.10433 (9)	0.32342(7)	0.0217 (3)
C16A	0.5755 0.57737(14)	0.1102 0.04462 (0)	0.3723 0.20348 (8)	0.020
	0.57757 (14)	0.04402(9)	0.23346 (6)	0.0243(3)
	0.0270 0.53324 (14)	0.0110	0.3138	$0.029^{\circ}$
	0.5542	-0.0016	0.22813 (8)	0.0230 (3)
	0.3342 0.45901 (14)	-0.0040	0.2040 0.10785 (7)	$0.028^{\circ}$
	0.43801 (14)	0.08320 (8)	0.19785 (7)	0.0219 (3)
HI8A C10A	0.4278 0.42724 (12)	0.0784	0.1541	$0.020^{*}$
CI9A UI0A	0.42/34 (13)	0.14551 (8)	0.23259(7)	0.0184 (2)
HI9A	0.3771	0.1791	0.2120	0.022*
C20A	0.55895 (12)	0.30004 (8)	0.37121 (7)	0.0171 (2)
C21A	0.56048 (13)	0.36082 (8)	0.41418 (7)	0.0208 (3)
H2IA	0.5007	0.3686	0.4332	0.025*
C22A	0.65110 (15)	0.40982 (9)	0.42867 (8)	0.0267 (3)
H22A	0.6512	0.4509	0.4567	0.032*
C23A	0.74138 (14)	0.39751 (10)	0.40137 (9)	0.0288 (3)
H23A	0.8021	0.4302	0.4112	0.035*
C24A	0.74112 (14)	0.33647 (10)	0.35934 (9)	0.0271 (3)
H24A	0.8021	0.3279	0.3415	0.033*
C25A	0.64965 (13)	0.28783 (9)	0.34373 (8)	0.0216 (3)
H25A	0.6492	0.2474	0.3151	0.026*
C26A	0.13545 (13)	0.13752 (9)	0.67551 (7)	0.0226 (3)
H26A	0.2071	0.1510	0.6666	0.027*
C27A	0.1157 (2)	0.18750 (15)	0.73067 (11)	0.0520 (5)
H27A	0.1225	0.2381	0.7213	0.078*
H27B	0.1698	0.1780	0.7689	0.078*
H27C	0.0424	0.1781	0.7368	0.078*
C28A	0.13514 (19)	0.05692 (11)	0.68819 (11)	0.0395 (5)
H28A	0.1436	0.0279	0.6509	0.059*
H28B	0.0660	0.0437	0.6982	0.059*
H28C	0.1956	0.0477	0.7235	0.059*
C29A	-0.16430 (12)	0.16323 (9)	0.53451 (8)	0.0219 (3)

H29A	-0.1631	0.1786	0.5791	0.026*
C30A	-0.17230 (18)	0.08025 (11)	0.52515 (13)	0.0427 (4)
H30A	-0.1106	0.0593	0.5537	0.064*
H30B	-0.1707	0.0654	0.4818	0.064*
H30C	-0.2407	0.0632	0.5340	0.064*
C31A	-0.25866 (18)	0.19967 (14)	0.49048 (12)	0.0440 (4)
H31A	-0.2528	0.2521	0.4996	0.066*
H31B	-0.3283	0.1818	0.4969	0.066*
H31C	-0.2549	0.1883	0.4469	0.066*
C32A	0.04283 (14)	0.35253 (8)	0.59024 (8)	0.0223 (3)
H32A	0.0905	0.3536	0.5596	0.027*
C33A	-0.07353 (16)	0.37404 (10)	0.55800 (10)	0.0302 (4)
H33A	-0.0996	0.3430	0.5197	0.045*
H33B	-0.0728	0.4246	0.5474	0.045*
H33C	-0.1221	0.3683	0.5865	0.045*
C34A	0.09399 (19)	0.40151 (10)	0.64906 (10)	0.0348 (4)
H34A	0.1656	0.3833	0.6688	0.052*
H34B	0.0466	0.4014	0.6787	0.052*
H34C	0.1023	0.4509	0.6370	0.052*
C35A	0.05387 (13)	0.25422 (8)	0.38600 (7)	0.0198 (3)
C36A	0.09066 (13)	0.38836 (8)	0.44165 (7)	0.0186 (2)
C37A	0.30942 (13)	0.36302 (8)	0.48535 (7)	0.0194 (2)
C38A	0.44768 (12)	0.23894 (8)	0.50448 (7)	0.0181 (2)
C39A	0.40018 (13)	0.09858 (9)	0.44396 (7)	0.0209 (3)
C40A	0.22252 (12)	0.15734 (8)	0.36194 (7)	0.0189 (2)
C41A	0.10377 (13)	0.12908 (8)	0.45433 (7)	0.0207 (3)
C42A	0.27394 (13)	0.11760 (9)	0.55825 (7)	0.0208 (3)
C43A	0.28449 (13)	0.26821 (9)	0.58020 (7)	0.0211 (3)
Ru1B	0.339824 (9)	0.737543 (6)	0.020451 (5)	0.01278 (2)
Ru2B	0.554997 (9)	0.807238 (6)	0.065368 (5)	0.01303 (2)
Ru3B	0.409502 (9)	0.797940 (6)	0.148373 (5)	0.01333 (2)
As1B	0.354067 (12)	0.725193 (8)	-0.090391 (7)	0.01419 (2)
As2B	0.617117 (12)	0.757903 (8)	-0.027337 (7)	0.01422 (2)
P1B	0.24434 (3)	0.79901 (2)	0.179505 (17)	0.01526 (6)
O1B	0.48228 (10)	0.60037 (6)	0.04909 (6)	0.0230 (2)
O2B	0.12319 (10)	0.65451 (7)	0.01519 (7)	0.0297 (3)
O3B	0.19725 (10)	0.87497 (6)	-0.00891 (6)	0.0250 (2)
O4B	0.43815 (10)	0.92849 (6)	-0.01740 (6)	0.0259 (2)
O5B	0.72668 (10)	0.92191 (7)	0.13537 (6)	0.0285 (2)
O6B	0.67389 (10)	0.68923 (6)	0.15124 (6)	0.0234 (2)
O7B	0.44806 (12)	0.63705 (7)	0.18197 (6)	0.0296 (3)
O8B	0.58619 (11)	0.84901 (8)	0.26636 (6)	0.0324 (3)
O9B	0.35312 (11)	0.95402 (6)	0.10515 (6)	0.0254 (2)
O10B	0.24208 (10)	0.83223 (7)	0.24973 (6)	0.0260 (2)
O11B	0.17395 (9)	0.72522 (6)	0.17894 (6)	0.0217 (2)
O12B	0.15674 (9)	0.84427 (6)	0.13157 (5)	0.0203 (2)
C1B	0.33933 (12)	0.62820 (7)	-0.13510 (7)	0.0163 (2)
C2B	0.35030 (13)	0.62110 (8)	-0.19829 (7)	0.0201 (3)
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H2BA	0.3574	0.6630	-0.2200	0.024*
C3B	0.35059 (14)	0.55150 (9)	-0.22884 (7)	0.0227 (3)
H3BA	0.3578	0.5468	-0.2709	0.027*
C4B	0.34000 (13)	0.48894 (9)	-0.19637 (8)	0.0227 (3)
H4BA	0.3419	0.4423	-0.2165	0.027*
C5B	0.32664 (12)	0.49591 (8)	-0.13416(7)	0.0200 (3)
H5BA	0.3180	0.4540	-0.1129	0.024*
C6B	0.32616 (12)	0.56564 (8)	-0.10337(7)	0.0177 (2)
H6BA	0.3170	0.5702	-0.0616	0.021*
C7B	0.25050 (12)	0.78285 (8)	-0.14899(7)	0.0169 (2)
C8B	0.27299 (15)	0.85673 (9)	-0.15500(9)	0.0279 (3)
H8BA	0.3416	0.8778	-0.1343	0.033*
C9B	0.19267 (16)	0.89929 (9)	-0.19204(9)	0.0317(4)
H9BA	0.2081	0.9487	-0.1961	0.0317 (1)
C10B	0.09035(15)	0.86873 (9)	-0.22268(8)	0.020
H10B	0.0372	0.8077	-0.2477	0.0272 (3)
C11B	0.0572 0.06720 (13)	0.79510 (9)	-0.21592(8)	0.033
H11B	-0.00720(13)	0.79510 (9)	-0.2360	0.0235 (3)
C12B	0.0020 0.14688 (12)	0.7743 0.75218 (8)	-0.17030(7)	0.028
U12D	0.14088 (12)	0.75218 (8)	-0.1750	0.0197(3)
П12В	0.1310 0.40909 (12)	0.7029	-0.1730	$0.024^{\circ}$
	0.49808 (12)	0.73611 (8)	-0.10303 (7)	0.0188 (2)
HISC	0.5175	0.7258	-0.1382	0.023*
HI3D	0.4912	0.80/6	-0.1199	0.023*
CI4B	0.6/501 (13)	0.65916 (8)	-0.028/9(/)	0.0190 (3)
CISB	0.7/084 (15)	0.64590 (9)	0.01763 (8)	0.0255 (3)
HI5B	0.7999	0.6823	0.0495	0.031*
C16B	0.82291 (18)	0.57824 (10)	0.01631 (10)	0.0337 (4)
H16B	0.8874	0.5696	0.0468	0.040*
C17B	0.77798 (18)	0.52375 (10)	-0.03081 (11)	0.0370 (5)
H17B	0.8129	0.4787	-0.0320	0.044*
C18B	0.68180 (17)	0.53615 (10)	-0.07589 (11)	0.0357 (4)
H18B	0.6513	0.4990	-0.1067	0.043*
C19B	0.63018 (14)	0.60414 (9)	-0.07547 (9)	0.0265 (3)
H19B	0.5660	0.6126	-0.1063	0.032*
C20B	0.73519 (12)	0.81087 (8)	-0.05346 (7)	0.0173 (2)
C21B	0.80847 (14)	0.77348 (9)	-0.08396 (8)	0.0247 (3)
H21B	0.7994	0.7229	-0.0937	0.030*
C22B	0.89520 (16)	0.81160 (10)	-0.09990 (10)	0.0316 (4)
H22B	0.9443	0.7864	-0.1201	0.038*
C23B	0.90882 (15)	0.88701 (10)	-0.08582 (10)	0.0307 (4)
H23B	0.9672	0.9124	-0.0963	0.037*
C24B	0.83519 (14)	0.92459 (9)	-0.05607 (9)	0.0267 (3)
H24B	0.8437	0.9753	-0.0472	0.032*
C25B	0.74863 (12)	0.88669 (8)	-0.03947 (8)	0.0208 (3)
H25B	0.6999	0.9120	-0.0191	0.025*
C26B	0.31227 (16)	0.89327 (10)	0.28401 (9)	0.0305 (4)
H26B	0.3632	0.9091	0.2583	0.037*
C27B	0.2382 (2)	0.95604 (15)	0.29488 (11)	0.0520 (5)

H27D	0.1988	0.9722	0.2544	0.078*
H27E	0.2828	0.9962	0.3190	0.078*
H27F	0.1858	0.9397	0.3182	0.078*
C28B	0.37867 (18)	0.86573 (14)	0.34603 (9)	0.0392 (4)
H28D	0.4231	0.8251	0.3369	0.059*
H28E	0.3287	0.8498	0.3710	0.059*
H28F	0.4263	0.9048	0.3696	0.059*
C29B	0.20585 (15)	0.66959 (10)	0.22553 (10)	0.0293 (4)
H29B	0.2850	0.6757	0.2464	0.035*
C30B	0.18501 (18)	0.59571 (11)	0.18843 (13)	0.0427 (4)
H30D	0.2279	0.5930	0.1564	0.064*
H30E	0.1074	0.5900	0.1682	0.064*
H30F	0.2065	0.5571	0.2170	0.064*
C31B	0.13696 (18)	0.67900 (14)	0.27457 (12)	0.0440 (4)
H31D	0.1547	0.7259	0.2980	0.066*
H31E	0.1527	0.6403	0.3035	0.066*
H31F	0.0594	0.6769	0.2535	0.066*
C32B	0.04202 (12)	0.85458 (9)	0.13787 (8)	0.0227 (3)
H32B	0.0375	0.8471	0.1819	0.027*
C33B	-0.03334 (15)	0.79981 (11)	0.09303 (11)	0.0365 (4)
H33D	-0.0127	0.7507	0.1045	0.055*
H33E	-0.0266	0.8058	0.0501	0.055*
H33F	-0.1089	0.8078	0.0958	0.055*
C34B	0.01410 (15)	0.93339 (10)	0.12292 (11)	0.0319 (4)
H34D	0.0661	0.9663	0.1518	0.048*
H34E	-0.0599	0.9433	0.1278	0.048*
H34F	0.0184	0.9407	0.0797	0.048*
C35B	0.43313 (12)	0.65362 (8)	0.04108 (7)	0.0171 (2)
C36B	0.20505 (12)	0.68588 (8)	0.01746 (7)	0.0191 (2)
C37B	0.25441 (12)	0.82674 (8)	0.00457 (7)	0.0179 (2)
C38B	0.47542 (12)	0.88203 (8)	0.01420 (7)	0.0184 (2)
C39B	0.66404 (12)	0.87716 (8)	0.10905 (7)	0.0182 (2)
C40B	0.62378 (12)	0.73118 (8)	0.11932 (7)	0.0178 (2)
C41B	0.43382 (13)	0.69593 (8)	0.16652 (7)	0.0203 (3)
C42B	0.51685 (13)	0.83132 (9)	0.22284 (7)	0.0213 (3)
C43B	0.37617 (12)	0.89545 (8)	0.11856 (7)	0.0188 (2)

Atomic displacement parameters  $(Å^2)$ 

	$U^{11}$	$U^{22}$	$U^{33}$	$U^{12}$	$U^{13}$	$U^{23}$
Ru1A	0.01498 (4)	0.01348 (4)	0.01328 (4)	0.00320 (3)	0.00426 (3)	0.00231 (3)
Ru2A	0.01373 (4)	0.01421 (4)	0.01253 (4)	0.00283 (3)	0.00234 (3)	0.00207 (3)
Ru3A	0.01465 (4)	0.01462 (4)	0.01440 (4)	0.00116 (3)	0.00371 (4)	0.00334 (3)
As1A	0.01435 (6)	0.01388 (6)	0.01247 (6)	0.00331 (4)	0.00323 (4)	0.00196 (4)
As2A	0.01417 (6)	0.01534 (6)	0.01208 (6)	0.00351 (4)	0.00263 (4)	0.00179 (4)
P1A	0.01446 (14)	0.01624 (15)	0.01473 (14)	0.00120 (11)	0.00285 (12)	0.00451 (11)
O1A	0.0223 (5)	0.0260 (6)	0.0318 (6)	0.0010 (4)	0.0012 (5)	0.0003 (5)
O2A	0.0316 (6)	0.0217 (5)	0.0361 (7)	0.0102 (5)	0.0170 (5)	0.0054 (5)

# supporting information

O3A	0.0284 (6)	0.0238 (5)	0.0243 (6)	0.0000 (4)	-0.0017 (5)	0.0002 (4)
O4A	0.0218 (5)	0.0287 (6)	0.0209 (5)	0.0021 (4)	-0.0001 (4)	-0.0016 (4)
O5A	0.0432 (8)	0.0236 (6)	0.0329 (7)	0.0138 (5)	0.0039 (6)	0.0060 (5)
06A	0.0214 (5)	0.0308 (6)	0.0246 (6)	0.0032 (4)	-0.0008 (4)	-0.0047 (5)
O7A	0.0334 (6)	0.0240 (6)	0.0284 (6)	-0.0051(5)	0.0038 (5)	-0.0012(5)
O8A	0.0347 (7)	0.0281 (6)	0.0285 (6)	0.0116 (5)	0.0059 (5)	0.0110 (5)
09A	0.0289 (6)	0.0367 (7)	0.0241 (6)	-0.0104 (5)	0.0054 (5)	-0.0039 (5)
010A	0.0167 (4)	0.0221 (5)	0.0202 (5)	0.0008 (4)	0.0026 (4)	0.0104 (4)
011A	0.0148 (4)	0.0246 (5)	0.0178 (5)	0.0010 (4)	0.0020 (4)	0.0049 (4)
012A	0.0248 (5)	0.0181 (5)	0.0181 (5)	0.0012 (4)	0.0084 (4)	0.0034 (4)
C1A	0.0148 (5)	0.0143 (5)	0.0159 (6)	0.0017 (4)	0.0011 (4)	0.0023 (4)
C2A	0.0174 (6)	0.0307 (8)	0.0205 (7)	0.0070 (5)	0.0032 (5)	0.0026 (6)
C3A	0.0183 (7)	0.0320 (8)	0.0293 (8)	0.0089 (6)	0.0017 (6)	0.0040 (6)
C4A	0.0165 (6)	0.0227 (7)	0.0267 (7)	0.0005 (5)	-0.0026(5)	0.0089 (6)
C5A	0.0242 (7)	0.0336 (8)	0.0186 (7)	0.0049 (6)	0.0014 (6)	0.0080 (6)
C6A	0.0221 (7)	0.0304 (8)	0.0178 (6)	0.0097 (6)	0.0034 (5)	0.0050 (5)
C7A	0.0160(5)	0.0167 (6)	0.0148 (5)	0.0018 (4)	0.0029 (4)	0.0031 (4)
C8A	0.0269 (7)	0.0210 (6)	0.0213 (7)	0.0017 (5)	0.0111 (6)	0.0025 (5)
C9A	0.0245 (7)	0.0235 (7)	0.0267(7)	-0.0003(5)	0.0113 (6)	0.0063 (6)
C10A	0.0216 (7)	0.0209 (6)	0.0230(7)	-0.0033(5)	0.0022 (5)	0.0034(5)
C11A	0.0265(7)	0.0208 (7)	0.0216 (7)	-0.0037(5)	0.0058 (6)	-0.0023(5)
C12A	0.0214 (6)	0.0192 (6)	0.0186 (6)	-0.0009(5)	0.0064 (5)	0.0003 (5)
C13A	0.0175 (6)	0.0171 (6)	0.0138 (5)	0.0054 (4)	0.0032 (4)	0.0018 (4)
C14A	0.0181 (6)	0.0155 (5)	0.0159 (6)	0.0027 (4)	0.0051(5)	0.0016 (4)
C15A	0.0234 (7)	0.0232 (7)	0.0178 (6)	0.0089 (5)	0.0031 (5)	0.0012 (5)
C16A	0.0272 (7)	0.0216 (7)	0.0243 (7)	0.0100 (6)	0.0057 (6)	0.0032 (5)
C17A	0.0314 (8)	0.0162 (6)	0.0230(7)	0.0043 (5)	0.0095 (6)	0.0011(5)
C18A	0.0324 (8)	0.0161 (6)	0.0172 (6)	0.0016 (5)	0.0056 (6)	0.0005 (5)
C19A	0.0240 (7)	0.0151 (6)	0.0158 (6)	0.0023 (5)	0.0033 (5)	0.0020 (4)
C20A	0.0158 (6)	0.0189 (6)	0.0159 (6)	0.0016 (4)	0.0013 (5)	0.0044 (5)
C21A	0.0211 (6)	0.0218 (6)	0.0188 (6)	0.0008 (5)	0.0029 (5)	0.0016 (5)
C22A	0.0288 (8)	0.0247 (7)	0.0233 (7)	-0.0051 (6)	-0.0007 (6)	0.0015 (6)
C23A	0.0226 (7)	0.0323 (8)	0.0292 (8)	-0.0071 (6)	-0.0008 (6)	0.0108 (7)
C24A	0.0178 (6)	0.0352 (8)	0.0294 (8)	0.0010 (6)	0.0050 (6)	0.0116 (7)
C25A	0.0185 (6)	0.0260 (7)	0.0215 (7)	0.0035 (5)	0.0059 (5)	0.0059 (5)
C26A	0.0204 (6)	0.0299 (7)	0.0183 (6)	0.0050 (5)	0.0036 (5)	0.0106 (5)
C27A	0.0609 (11)	0.0565 (10)	0.0322 (7)	0.0290 (9)	-0.0004 (7)	-0.0087 (7)
C28A	0.0400 (10)	0.0346 (10)	0.0419 (11)	0.0058 (8)	0.0000 (9)	0.0227 (8)
C29A	0.0155 (6)	0.0260 (7)	0.0242 (7)	-0.0007(5)	0.0044 (5)	0.0035 (5)
C30A	0.0312 (7)	0.0245 (6)	0.0691 (11)	-0.0042(5)	0.0036 (7)	0.0074 (6)
C31A	0.0319 (7)	0.0525 (9)	0.0538 (9)	0.0109 (6)	0.0167 (7)	0.0258 (8)
C32A	0.0283 (7)	0.0176 (6)	0.0243 (7)	-0.0001(5)	0.0124 (6)	0.0028 (5)
C33A	0.0337 (9)	0.0229 (7)	0.0383 (9)	0.0087 (6)	0.0150 (7)	0.0102 (7)
C34A	0.0459 (11)	0.0261 (8)	0.0351 (9)	-0.0104 (7)	0.0176 (8)	-0.0073 (7)
C35A	0.0209 (6)	0.0179 (6)	0.0214 (6)	0.0045 (5)	0.0061 (5)	0.0019 (5)
C36A	0.0214 (6)	0.0178 (6)	0.0185 (6)	0.0023 (5)	0.0075 (5)	0.0039 (5)
C37A	0.0212 (6)	0.0198 (6)	0.0174 (6)	0.0044 (5)	0.0044 (5)	0.0041 (5)
C38A	0.0183 (6)	0.0195 (6)	0.0167 (6)	0.0036 (5)	0.0044 (5)	0.0023 (5)
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C39A	0.0220 (7)	0.0221 (6)	0.0180 (6)	0.0051 (5)	0.0031 (5)	0.0026 (5)
C40A	0.0165 (6)	0.0206 (6)	0.0193 (6)	0.0043 (5)	0.0030 (5)	0.0022 (5)
C41A	0.0234 (7)	0.0190 (6)	0.0204 (6)	0.0017 (5)	0.0059 (5)	0.0038 (5)
C42A	0.0213 (6)	0.0222 (6)	0.0198 (6)	0.0021 (5)	0.0057 (5)	0.0049 (5)
C43A	0.0188 (6)	0.0249 (7)	0.0207 (6)	-0.0016 (5)	0.0064 (5)	0.0033 (5)
Ru1B	0.01209 (4)	0.01290 (4)	0.01270 (4)	-0.00062(3)	0.00130 (3)	0.00173 (3)
Ru2B	0.01213 (4)	0.01236 (4)	0.01426 (4)	-0.00093 (3)	0.00238 (3)	0.00084 (3)
Ru3B	0.01277 (4)	0.01367 (4)	0.01320 (4)	0.00093 (3)	0.00229 (3)	0.00055 (3)
As1B	0.01590 (6)	0.01336 (6)	0.01246 (6)	-0.00087(4)	0.00123 (5)	0.00213 (4)
As2B	0.01477 (6)	0.01318 (6)	0.01513 (6)	-0.00051 (4)	0.00402 (5)	0.00199 (4)
P1B	0.01455 (14)	0.01611 (15)	0.01531 (15)	0.00140 (11)	0.00363 (12)	0.00172 (11)
O1B	0.0216 (5)	0.0187 (5)	0.0261 (5)	0.0022 (4)	-0.0001 (4)	0.0031 (4)
O2B	0.0191 (5)	0.0244 (6)	0.0455 (8)	-0.0045(4)	0.0067 (5)	0.0045 (5)
O3B	0.0235 (5)	0.0208 (5)	0.0285 (6)	0.0046 (4)	0.0012 (4)	0.0027 (4)
O4B	0.0250(5)	0.0199 (5)	0.0303 (6)	-0.0011(4)	-0.0002(5)	0.0078(4)
05B	0.0245 (6)	0.0249(6)	0.0321 (6)	-0.0076(4)	0.0002(0)	-0.0040(5)
06B	0.0236(5)	0.0220(5)	0.0229(5)	0 0024 (4)	0.0029(4)	0.0062(4)
07B	0.0230(3) 0.0389(7)	0.0220(5)	0.0235(6)	0.0021(1)	0.0191 (6)	0.0002(1)
08B	0.0233(6)	0.0230(3) 0.0414(7)	0.0266 (6)	0.0101(5)	-0.0048(5)	-0.0052(5)
09B	0.0295 (6)	0.0161(5)	0.0284 (6)	0.0002(3)	0.0023(5)	0.0002(0)
010B	0.0299(0) 0.0249(5)	0.0101(5) 0.0348(6)	0.0201(0)	0.0023(1) 0.0018(5)	0.0025(3)	-0.0049(4)
010B	0.0215(5) 0.0185(5)	0.0195(5)	0.0171(5) 0.0275(5)	-0.0011(4)	0.0050 (4)	0.0069 (4)
012B	0.0102(3) 0.0142(4)	0.0243(5)	0.0278(5)	0.0038(4)	0.0053(4)	0.00094(4)
C1B	0.0165 (6)	0.0215(5)	0.0157 (6)	0.0000(1)	0.00000(1)	0.00012(4)
C2B	0.0109(0) 0.0240(7)	0.0197(6)	0.0164 (6)	0.0001(1) 0.0009(5)	0.0018(5)	0.0012(1) 0.0020(5)
C3B	0.0210(7) 0.0257(7)	0.0237(7)	0.0180 (6)	0.0009(5)	0.0030(5)	-0.0020(5)
C4B	0.0237(7) 0.0213(7)	0.0237(7)	0.0259(7)	0.0053(5)	0.0020 (6)	-0.0023(5)
C5B	0.0219(7)	0.0154 (6)	0.0237(7)	0.0023(5)	0.0020(0)	0.0025(5)
C6B	0.0174 (6)	0.0169 (6)	0.0174(6)	0.0021(5)	0.0009(5)	0.0025(5)
C7B	0.0195 (6)	0.0164 (6)	0.0134(5)	0.0002(0)	0.0003(0)	0.0020(0)
C8B	0.0190(8)	0.0193(7)	0.019 (8)	-0.0057(6)	-0.0082(6)	0.0070 (6)
C9B	0.0200(0) 0.0342(9)	0.0175(7)	0.0299(0) 0.0358(9)	-0.0037(6)	-0.0002(0)	0.0093 (6)
C10B	0.0312(9) 0.0279(8)	0.0227(7)	0.0266 (8)	0.0023(6)	-0.0098(6)	0.0082 (6)
C11B	0.0279 (6)	0.0227(7) 0.0243(7)	0.0200(0) 0.0242(7)	-0.0023(0)	-0.0007(5)	0.0065 (6)
C12B	0.0195 (6)	0.0213(7)	0.0212(7)	-0.0017(5)	0.0007(3)	0.0000(0)
C13B	0.0193 (6)	0.0177(0)	0.0154 (6)	-0.0017(5)	0.0019(5)	0.0031(5)
C14B	0.0218 (6)	0.0210(0)	0.0134(0) 0.0236(7)	0.003 + (3)	0.0025(5)	0.0035(5)
C15B	0.0210(0) 0.0304(8)	0.0149(0)	0.0250(7)	0.0069 (6)	0.0103(5)	0.0035(5)
C16B	0.0304(0)	0.0227(7)	0.0230(7) 0.0374(10)	0.0009(0)	0.0085 (8)	0.0074(0)
C17B	0.0402(10) 0.0456(11)	0.0287(3) 0.0184(7)	0.0571(13)	0.0149(7)	0.0100(0)	0.0103(7)
C17B	0.0430(11) 0.0372(10)	0.0134(7)	0.0571(13)	-0.0025(6)	0.0303(10)	-0.0054(7)
C10B	0.0372(10) 0.0231(7)	0.0170(7)	0.0380(13)	-0.0025(0)	0.0233(9)	-0.0034(7)
C20B	0.0231(7) 0.0158(6)	0.0199(7) 0.0187(6)	0.0380(9)	-0.0025(3)	0.0119(7)	0.0043(0)
C20B	0.0138(0) 0.0277(7)	0.0107(0) 0.0212(7)	0.0101(0) 0.0207(8)	0.0000 (4)	0.0071(3)	0.0033(3)
C21D	0.0277(7)	0.0213(7)	0.027(0) 0.0442(10)	0.0011(0) 0.0042(6)	0.0132(0) 0.0227(9)	0.00+2(0)
C22D	0.0303(0)	0.0203 (0)	0.0443(10)	0.0042(0)	0.0237(0)	0.0099(7)
C23D	0.0230(7)	0.0292(6) 0.0106(7)	0.0400(10)	-0.0018(0)	0.0100(7)	0.0103(7)
C24D	0.0223(7)	0.0190(7)	0.0400 (9)	-0.0012(3)	0.0099(7)	0.0102(0)
UZ3B	0.0164 (0)	0.01/1(0)	0.0283 (7)	0.0000 (3)	0.0071(3)	0.0009 (3)

C26B	0.0317 (9)	0.0329 (9)	0.0235 (8)	0.0074 (7)	0.0020 (6)	-0.0100 (6)
C27B	0.0609 (11)	0.0565 (10)	0.0322 (7)	0.0290 (9)	-0.0004 (7)	-0.0087 (7)
C28B	0.0327 (10)	0.0588 (13)	0.0236 (8)	0.0047 (9)	0.0029 (7)	-0.0031 (8)
C29B	0.0235 (7)	0.0282 (8)	0.0414 (10)	0.0070 (6)	0.0139 (7)	0.0184 (7)
C30B	0.0312 (7)	0.0245 (6)	0.0691 (11)	-0.0042 (5)	0.0036 (7)	0.0074 (6)
C31B	0.0319 (7)	0.0525 (9)	0.0538 (9)	0.0109 (6)	0.0167 (7)	0.0258 (8)
C32B	0.0143 (6)	0.0250 (7)	0.0309 (8)	0.0024 (5)	0.0075 (5)	0.0096 (6)
C33B	0.0189 (7)	0.0316 (9)	0.0541 (12)	0.0000 (6)	-0.0023 (8)	0.0058 (8)
C34B	0.0204 (7)	0.0247 (8)	0.0514 (11)	0.0059 (6)	0.0074 (7)	0.0113 (7)
C35B	0.0159 (6)	0.0191 (6)	0.0153 (6)	-0.0015 (5)	0.0011 (5)	0.0025 (4)
C36B	0.0184 (6)	0.0159 (6)	0.0221 (6)	0.0010 (5)	0.0026 (5)	0.0026 (5)
C37B	0.0170 (6)	0.0174 (6)	0.0178 (6)	-0.0021 (5)	0.0011 (5)	0.0009 (5)
C38B	0.0156 (6)	0.0165 (6)	0.0220 (6)	-0.0019 (4)	0.0022 (5)	0.0006 (5)
C39B	0.0168 (6)	0.0179 (6)	0.0192 (6)	0.0001 (5)	0.0028 (5)	0.0008 (5)
C40B	0.0174 (6)	0.0172 (6)	0.0184 (6)	-0.0023 (5)	0.0038 (5)	0.0001 (5)
C41B	0.0205 (6)	0.0221 (6)	0.0208 (6)	0.0043 (5)	0.0094 (5)	0.0042 (5)
C42B	0.0197 (6)	0.0231 (7)	0.0203 (6)	0.0023 (5)	0.0034 (5)	0.0001 (5)
C43B	0.0190 (6)	0.0186 (6)	0.0173 (6)	-0.0003 (5)	0.0016 (5)	-0.0005 (5)

## Geometric parameters (Å, °)

Ru1A—C36A	1.8731 (14)	Ru1B—C36B	1.8914 (15)
Ru1A—C37A	1.9246 (16)	Ru1B—C35B	1.9292 (14)
Ru1A—C35A	1.9386 (16)	Ru1B—C37B	1.9445 (14)
Ru1A—As1A	2.4596 (2)	Ru1B—As1B	2.4297 (2)
Ru1A—Ru2A	2.8781 (2)	Ru1B—Ru3B	2.8397 (2)
Ru1A—Ru3A	2.9093 (2)	Ru1B—Ru2B	2.8685 (2)
Ru2A—C39A	1.8945 (15)	Ru2B—C39B	1.8876 (15)
Ru2A—C40A	1.9225 (15)	Ru2B—C38B	1.9315 (15)
Ru2A—C38A	1.9343 (15)	Ru2B—C40B	1.9340 (15)
Ru2A—As2A	2.4136 (2)	Ru2B—As2B	2.4223 (2)
Ru2A—Ru3A	2.8106 (2)	Ru2B—Ru3B	2.8200 (2)
Ru3A—C42A	1.8891 (15)	Ru3B—C42B	1.8955 (16)
Ru3A—C43A	1.9298 (16)	Ru3B—C41B	1.9287 (15)
Ru3A—C41A	1.9491 (16)	Ru3B—C43B	1.9359 (15)
Ru3A—P1A	2.2926 (4)	Ru3B—P1B	2.2877 (4)
As1A—C7A	1.9494 (14)	As1B—C7B	1.9404 (14)
As1A—C1A	1.9532 (13)	As1B—C1B	1.9426 (14)
As1A—C13A	1.9638 (13)	As1B—C13B	1.9711 (14)
As2A—C14A	1.9363 (14)	As2B—C14B	1.9369 (14)
As2A—C20A	1.9426 (14)	As2B—C20B	1.9470 (13)
As2A—C13A	1.9626 (13)	As2B—C13B	1.9583 (15)
P1A—O11A	1.5839 (11)	P1B-012B	1.5860 (11)
P1A—O10A	1.5935 (10)	P1B-011B	1.5909 (11)
P1A—O12A	1.6003 (11)	P1B	1.5997 (12)
01A—C35A	1.142 (2)	O1B—C35B	1.1464 (18)
O2A—C36A	1.1505 (18)	O2B—C36B	1.1469 (18)
O3A—C37A	1.149 (2)	O3B—C37B	1.1384 (18)

O4A—C38A	1.1437 (19)	O4B—C38B	1.1486 (18)
O5A—C39A	1.1441 (19)	O5B—C39B	1.1470 (18)
O6A—C40A	1.1489 (19)	O6B—C40B	1.1452 (18)
O7A—C41A	1.136 (2)	O7B—C41B	1.1426 (18)
08A—C42A	1.1410 (19)	O8B—C42B	1.143 (2)
09A - C43A	1 145 (2)	O9B-C43B	1 1406 (18)
010A - C26A	1.113(2) 1 4584 (19)	O10B-C26B	1.1100(10) 1456(2)
0114 - C294	1.4697 (18)	O11B-C29B	1.150(2) 1.4596(19)
O12A C22A	1.4097(10) 1.4505(18)	012P C22P	1.4590(19)
C1A = C6A	1.4393(10) 1.200(2)	C1R C6R	1.4038(18) 1.3887(10)
C1A = C0A	1.390(2) 1.201(2)	CID-COB	1.306/(19)
CIA—CZA	1.391 (2)	C1D - C2D	1.390 (2)
C2A—C3A	1.396 (2)		1.390 (2)
C2A—H2AA	0.9300	C2B—H2BA	0.9300
C3A—C4A	1.383 (2)	C3B—C4B	1.391 (2)
СЗА—НЗАА	0.9300	СЗВ—НЗВА	0.9300
C4A—C5A	1.379 (2)	C4B—C5B	1.385 (2)
C4A—H4AA	0.9300	C4B—H4BA	0.9300
C5A—C6A	1.397 (2)	C5B—C6B	1.395 (2)
С5А—Н5АА	0.9300	C5B—H5BA	0.9300
С6А—Н6АА	0.9300	C6B—H6BA	0.9300
C7A—C8A	1.392 (2)	C7B—C8B	1.389 (2)
C7A—C12A	1.395 (2)	C7B—C12B	1.392 (2)
C8A—C9A	1.392 (2)	C8B—C9B	1.393 (2)
С8А—Н8АА	0.9300	C8B—H8BA	0.9300
C9A—C10A	1.383 (2)	C9B—C10B	1.380(2)
С9А—Н9АА	0.9300	C9B—H9BA	0.9300
C10A - C11A	1 391 (2)	C10B-C11B	1389(2)
C10A - H10A	0.9300	C10B—H10B	0.9300
	1.301(2)		1.389(2)
$C_{11A} = H_{11A}$	0.0300		1.307(2)
	0.9300		0.9300
C12A $H12A$	0.9300	C12D—H12D	0.9300
	0.9700		0.9700
CI3A—HI3B	0.9700	CI3B—HI3D	0.9700
C14A—C19A	1.391 (2)	C14B—C19B	1.390 (2)
C14A—C15A	1.397 (2)	C14B—C15B	1.396 (2)
C15A—C16A	1.388 (2)	C15B—C16B	1.393 (2)
C15A—H15A	0.9300	C15B—H15B	0.9300
C16A—C17A	1.387 (2)	C16B—C17B	1.389 (3)
C16A—H16A	0.9300	C16B—H16B	0.9300
C17A—C18A	1.385 (2)	C17B—C18B	1.380 (3)
C17A—H17A	0.9300	C17B—H17B	0.9300
C18A—C19A	1.393 (2)	C18B—C19B	1.394 (2)
C18A—H18A	0.9300	C18B—H18B	0.9300
C19A—H19A	0.9300	C19B—H19B	0.9300
C20A—C21A	1.394 (2)	C20B—C21B	1.392 (2)
C20A—C25A	1.394 (2)	C20B—C25B	1.393 (2)
C21A—C22A	1.391 (2)	C21B—C22B	1.390 (2)
C21A—H21A	0.9300	C21B—H21B	0.9300

C22A—C23A	1.387 (3)	C22B—C23B	1.386 (3)
C22A—H22A	0.9300	C22B—H22B	0.9300
C23A—C24A	1.388 (3)	C23B—C24B	1.387 (2)
C23A—H23A	0.9300	C23B—H23B	0.9300
C24A—C25A	1.395 (2)	C24B—C25B	1.393 (2)
C24A—H24A	0.9300	C24B—H24B	0.9300
C25A—H25A	0.9300	C25B—H25B	0.9300
C26A—C28A	1.506 (2)	C26B—C27B	1.507 (3)
C26A—C27A	1.516 (3)	C26B—C28B	1.514 (3)
C26A—H26A	0.9800	C26B—H26B	0.9800
C27A—H27A	0.9600	C27B—H27D	0.9600
C27A—H27B	0.9600	C27B—H27E	0.9600
C27A—H27C	0.9600	C27B—H27F	0.9600
C28A—H28A	0.9600	C28B—H28D	0.9600
C28A—H28B	0.9600	C28B—H28E	0.9600
C28A—H28C	0.9600	C28B—H28F	0.9600
C29A—C31A	1.509 (3)	C29B—C31B	1.503 (3)
C29A—C30A	1.509 (3)	C29B—C30B	1.509 (3)
C29A—H29A	0.9800	C29B—H29B	0.9800
C30A—H30A	0.9600	C30B—H30D	0.9600
C30A—H30B	0.9600	C30B—H30E	0.9600
C30A—H30C	0.9600	C30B—H30F	0.9600
C31A—H31A	0.9600	C31B—H31D	0.9600
C31A—H31B	0.9600	C31B—H31E	0.9600
C31A—H31C	0.9600	C31B—H31F	0.9600
C32A—C33A	1.511 (3)	C32B—C33B	1.499 (3)
C32A—C34A	1.511 (3)	C32B—C34B	1.512 (2)
C32A—H32A	0.9800	C32B—H32B	0.9800
С33А—Н33А	0.9600	C33B—H33D	0.9600
С33А—Н33В	0.9600	C33B—H33E	0.9600
С33А—Н33С	0.9600	C33B—H33F	0.9600
C34A—H34A	0.9600	C34B—H34D	0.9600
C34A—H34B	0.9600	C34B—H34E	0.9600
C34A—H34C	0.9600	C34B—H34F	0.9600
C36A—Ru1A—C37A	90.69 (6)	C36B—Ru1B—C35B	95.01 (6)
C36A—Ru1A—C35A	87.60 (6)	C36B—Ru1B—C37B	88.31 (6)
C37A—Ru1A—C35A	167.69 (6)	C35B—Ru1B—C37B	175.52 (6)
C36A—Ru1A—As1A	98.87 (4)	C36B—Ru1B—As1B	102.23 (5)
C37A—Ru1A—As1A	93.49 (4)	C35B—Ru1B—As1B	91.41 (4)
C35A—Ru1A—As1A	98.82 (4)	C37B—Ru1B—As1B	90.82 (4)
C36A—Ru1A—Ru2A	168.90 (4)	C36B—Ru1B—Ru3B	105.09 (5)
C37A—Ru1A—Ru2A	83.29 (4)	C35B—Ru1B—Ru3B	90.48 (4)
C35A—Ru1A—Ru2A	96.28 (4)	C37B—Ru1B—Ru3B	85.74 (4)
As1A—Ru1A—Ru2A	90.831 (5)	As1B—Ru1B—Ru3B	152.336 (6)
C36A—Ru1A—Ru3A	112.69 (4)	C36B—Ru1B—Ru2B	162.49 (5)
C37A—Ru1A—Ru3A	89.28 (4)	C35B—Ru1B—Ru2B	78.46 (4)
C35A—Ru1A—Ru3A	80.16 (4)	C37B—Ru1B—Ru2B	97.50 (4)

As1A—Ru1A—Ru3A	148.290 (6)	As1B—Ru1B—Ru2B	94.217 (5)
Ru2A—Ru1A—Ru3A	58.107 (4)	Ru3B—Ru1B—Ru2B	59.208 (4)
C39A—Ru2A—C40A	92.15 (6)	C39B—Ru2B—C38B	91.23 (6)
C39A—Ru2A—C38A	93.44 (6)	C39B—Ru2B—C40B	91.03 (6)
C40A—Ru2A—C38A	173.78 (6)	C38B—Ru2B—C40B	175.25 (6)
C39A—Ru2A—As2A	100.95 (5)	C39B—Ru2B—As2B	106.85 (4)
C40A—Ru2A—As2A	85.40 (4)	C38B—Ru2B—As2B	89.22 (4)
C38A—Ru2A—As2A	96.18 (4)	C40B—Ru2B—As2B	94.12 (4)
C39A—Ru2A—Ru3A	100.87 (5)	C39B—Ru2B—Ru3B	103.35 (4)
C40A—Ru2A—Ru3A	92.12 (4)	C38B—Ru2B—Ru3B	96.62 (4)
C38A—Ru2A—Ru3A	84.17 (4)	C40B—Ru2B—Ru3B	78.79 (4)
As2A—Ru2A—Ru3A	158.123 (6)	As2B—Ru2B—Ru3B	149.091 (6)
C39A—Ru2A—Ru1A	159.78 (5)	C39B—Ru2B—Ru1B	158.69 (4)
C40A—Ru2A—Ru1A	79.49 (4)	C38B—Ru2B—Ru1B	78.77 (4)
C38A—Ru2A—Ru1A	94.34 (4)	C40B—Ru2B—Ru1B	97.72 (4)
As2A—Ru2A—Ru1A	96.728 (5)	As2B—Ru2B—Ru1B	91.922 (5)
Ru3A—Ru2A—Ru1A	61.503 (4)	Ru3B—Ru2B—Ru1B	59.889 (4)
C42A—Ru3A—C43A	91.93 (7)	C42B—Ru3B—C41B	92.30 (7)
C42A—Ru3A—C41A	90.06 (7)	C42B—Ru3B—C43B	94.97 (6)
C43A—Ru3A—C41A	177.11 (6)	C41B—Ru3B—C43B	172.34 (6)
C42A—Ru3A—P1A	99.67 (5)	C42B— $Ru3B$ — $P1B$	104.57 (5)
C43A—Ru3A—P1A	90.64 (4)	C41B—Ru3B—P1B	92.60 (4)
C41A—Ru3A—P1A	86.96 (5)	C43B—Ru3B—P1B	87.93 (4)
C42A—Ru3A—Ru2A	89.59 (4)	C42B—Ru3B—Ru2B	94.91 (5)
C43A—Ru3A—Ru2A	96.93 (4)	C41B—Ru3B—Ru2B	97.33 (4)
C41A—Ru3A—Ru2A	85.17 (4)	C43B—Ru3B—Ru2B	79.65 (4)
P1A—Ru3A—Ru2A	167.852 (10)	P1B—Ru3B—Ru2B	157.752 (11)
C42A—Ru3A—Ru1A	149.34 (5)	C42B—Ru3B—Ru1B	154.24 (5)
C43A—Ru3A—Ru1A	86.14 (5)	C41B—Ru3B—Ru1B	82.95 (5)
C41A—Ru3A—Ru1A	93.21 (4)	C43B—Ru3B—Ru1B	89.44 (4)
P1A—Ru3A—Ru1A	110.938 (10)	P1B—Ru3B—Ru1B	100.939 (10)
Ru2A—Ru3A—Ru1A	60.390 (4)	Ru2B—Ru3B—Ru1B	60.904 (4)
C7A—As1A—C1A	96.02 (6)	C7B—As1B—C1B	102.58 (6)
C7A—As1A—C13A	103.29 (6)	C7B—As1B—C13B	101.61 (6)
C1A—As1A—C13A	100.05 (6)	C1B—As1B—C13B	100.11 (6)
C7A—As1A—Ru1A	117.13 (4)	C7B—As1B—Ru1B	115.57 (4)
C1A—As1A—Ru1A	120.27 (4)	C1B—As1B—Ru1B	119.61 (4)
C13A—As1A—Ru1A	116.55 (4)	C13B—As1B—Ru1B	114.69 (4)
C14A—As2A—C20A	102.99 (6)	C14B—As2B—C20B	99.33 (6)
C14A—As2A—C13A	101.98 (6)	C14B—As2B—C13B	104.39(7)
C20A—As2A—C13A	102.67 (6)	C20B—As2B—C13B	100.91 (6)
C14A—As2A—Ru2A	114.16 (4)	C14B—As2B—Ru2B	119.98 (4)
C20A—As2A—Ru2A	122.44 (4)	C20B—As2B—Ru2B	118.26 (4)
C13A—As2A—Ru2A	110.20 (4)	C13B—As2B—Ru2B	111.40 (4)
O11A—P1A—O10A	100.18 (6)	O12B—P1B—O11B	100.20 (6)
O11A—P1A—O12A	107.60 (6)	O12B—P1B—O10B	106.00 (7)
O10A—P1A—O12A	98.03 (6)	O11B—P1B—O10B	98.48 (7)
O11A—P1A—Ru3A	110.48 (4)	O12B—P1B—Ru3B	109.20 (4)

O10A—P1A—Ru3A	117.33 (4)	O11B—P1B—Ru3B	121.52 (5)
O12A—P1A—Ru3A	120.75 (4)	O10B—P1B—Ru3B	118.99 (5)
C26A—O10A—P1A	123.23 (10)	C26B—O10B—P1B	124.91 (11)
C29A—O11A—P1A	125.30 (10)	C29B—O11B—P1B	123.45 (11)
C32A—O12A—P1A	123.41 (9)	C32B—O12B—P1B	122.81 (9)
C6A—C1A—C2A	119.26 (13)	C6B—C1B—C2B	119.84 (13)
C6A—C1A—As1A	121.58 (10)	C6B—C1B—As1B	120.55 (10)
C2A—C1A—As1A	118.78 (11)	C2B—C1B—As1B	119.44 (10)
C1A—C2A—C3A	119.99 (15)	C3B—C2B—C1B	120.03 (14)
C1A—C2A—H2AA	120.0	C3B—C2B—H2BA	120.0
СЗА—С2А—Н2АА	120.0	C1B—C2B—H2BA	120.0
C4A—C3A—C2A	120.36 (15)	C2B—C3B—C4B	119.92 (14)
С4А—С3А—НЗАА	119.8	C2B—C3B—H3BA	120.0
C2A - C3A - H3AA	119.8	C4B—C3B—H3BA	120.0
C5A - C4A - C3A	119.95 (14)	C5B-C4B-C3B	120.13 (14)
C5A - C4A - H4AA	120.0	C5B-C4B-H4BA	119.9
C3A - C4A - H4AA	120.0	C3B-C4B-H4BA	119.9
C4A - C5A - C6A	120.02 (15)	C4B-C5B-C6B	120 12 (14)
C4A - C5A - H5AA	120.02 (10)	C4B-C5B-H5BA	119.9
C6A - C5A - H5AA	120.0	C6B-C5B-H5BA	119.9
C1A - C6A - C5A	120.39 (14)	C1B-C6B-C5B	119.93 (13)
C1A - C6A - H6AA	119.8	C1B-C6B-H6BA	120.0
C5A - C6A - H6AA	119.8	C5B-C6B-H6BA	120.0
C8A - C7A - C12A	119.20 (13)	C8B-C7B-C12B	120.0 119.57(13)
C8A - C7A - As1A	122 40 (11)	C8B-C7B-As1B	120.75(11)
C12A - C7A - As1A	118 18 (10)	C12B - C7B - As1B	120.75(11) 119.22(10)
C9A - C8A - C7A	120.24(14)	C7B-C8B-C9B	119.22(10) 119.95(15)
C9A - C8A - H8AA	119.9	C7B-C8B-H8BA	120.0
C7A - C8A - H8AA	119.9	C9B-C8B-H8BA	120.0
C10A - C9A - C8A	120 54 (14)	C10B-C9B-C8B	120.0 120.52(15)
C10A - C9A - H9AA	110 7	C10B - C9B - H9BA	119.7
	119.7	C8B - C9B - H9BA	119.7
C9A - C10A - C11A	119.7	C9B-C10B-C11B	119.7
$C_{0A} = C_{10A} = H_{10A}$	120.3	$C^{0}B$ $C^{1}0B$ $H^{1}0B$	120.2
$C_{3A}$ $C_{10A}$ $H_{10A}$	120.3	$C_{11} C_{10} $	120.2
$C_{12A} = C_{10A} = M_{10A}$	120.3 120.20(14)	$C_{11}^{11}$ $C_{10}^{11}$ $C_{10}^{11}$ $C_{10}^{10}$ $C_{10}^{10}$	120.2 120.33(15)
$C_{12A} = C_{11A} = C_{10A}$	110.0	C12B $C11B$ $H11B$	120.33 (13)
$C_{12}A = C_{11}A = H_{11}A$	119.9	CIOR CIIR HIIR	119.8
$C_{11A} = C_{12A} = C_{7A}$	119.9	C11P C12P C7P	119.0
$C_{11A} = C_{12A} = C_{7A}$	120.23 (13)	$C_{11}^{11} = C_{12}^{12} = C_{13}^{12} = $	120.00 (13)
CTA = C12A = H12A	119.9	C7P $C12P$ $H12P$	120.0
C/A = C12A = H12A	119.9	C/D - C12D - D12D	120.0
$A_{32}A = C_{13}A = A_{31}A$	100.6	$A_{2}D = C_{1}D = H_{1}C$	111.95 (7)
$A_{s2A} = C_{13A} = H_{13A}$	109.0	$A_{s1D} = C_{12D} = H_{12C}$	109.2
$A_{2}A = C_{12}A = H_{12}B$	107.0	$A_{2}D = C_{12}D = U_{12}D$	109.2
$A_{2}A = C_{12}A = H_{12}B$	109.0	$A_{2}D = C_{12}D = H_{12}D$	109.2
H12A C12A H12D	109.0	ASID - CI3B - HI3D $H12C - C12D - H12D$	109.2
$\Pi I S A - U I S A - H I S B$	100.1	HI3U - HI3D	107.9
C19A—C14A—C15A	119.37 (13)	C19B—C14B—C15B	119.88 (14)

C19A-	C14AAs2A	123.08 (10)	C19B—C14B—As2B	123.26 (12)
C15A-	C14AAs2A	117.45 (11)	C15B—C14B—As2B	116.72 (12)
C16A-		120.40 (14)	C16B—C15B—C14B	120.07 (17)
C16A-		119.8	C16B—C15B—H15B	120.0
C14A-		119.8	C14B—C15B—H15B	120.0
C17A-		119.87 (14)	C17B—C16B—C15B	119.63 (19)
C17A-		120.1	C17B—C16B—H16B	120.2
C15A-		120.1	C15B—C16B—H16B	120.2
C18A-		120.07 (14)	C18B-C17B-C16B	120.38 (16)
C18A-		120.0	C18B—C17B—H17B	119.8
C16A-	-C17A-H17A	120.0	$C_{16B}$ $C_{17B}$ $H_{17B}$	119.8
C17A-	-C18A-C19A	120.28 (14)	C17B— $C18B$ — $C19B$	120 30 (19)
$C17A_{-}$	-C18A $-H18A$	119.9	C17B $C18B$ $H18B$	119.8
$C19A_{-}$	-C18A $-H18A$	119.9	C19B-C18B-H18B	119.8
$C14A_{-}$	-C19A-C18A	120.00(13)	C14B $C19B$ $C18B$	119.70 (18)
$C14A_{-}$	-C19A $-H19A$	120.00 (15)	C14B $C19B$ $H19B$	120.1
		120.0	C18B C19B H19B	120.1
$C_{10A}$	-C19A $-II19A$	120.0 110 66 (14)	$C_{10} = C_{10} = C$	120.1 110 71 (13)
C21A-	-C20A - C23A	119.00(14) 118.43(11)	$C_{21B} = C_{20B} = C_{23B}$	119.71(13) 120.78(11)
$C_{21A}$	-C20A - As2A	110.43(11) 121.01(11)	$C_{21D} = C_{20D} = A_{32D}$	120.78(11) 119.48(11)
$C_{23A}$	-C20A-AS2A	121.91(11) 120.19(15)	$C_{23} C_{20} $	119.48 (11)
$C_{22A}$	$-C_{21A}$ $-C_{20A}$	110.0	$C_{22} = C_{21} = C_{20} = C_{20}$	120.10 (13)
C22A-	$-C_{21A}$ $-H_{21A}$	119.9	$C_{22}D = C_{21}D = H_{21}D$	119.9
C20A-	$-C_{21}A$ $-H_{21}A$	119.9	$C_{20} = C_{21} = C$	119.9
C23A-	-C22A $-C21A$	120.07 (10)	$C_{23} = C_{22} = C_{21} = C_{21} = C_{21} = C_{22} = C$	120.21 (10)
C23A-	-C22A $H22A$	120.0	$C_{23}D = C_{22}D = H_{22}D$	119.9
C21A-	-C22A $-H22A$	120.0	$C_{21}D = C_{22}D = C_{24}D$	119.9
C24A-	-C23A-C22A	120.01 (16)	$C_{22}B = C_{23}B = C_{24}B$	119.85 (15)
C24A-	-C23A-H23A	120.0	C22B—C23B—H23B	120.1
C22A-	-C23A-H23A	120.0	$C_{24B}$ $C_{23B}$ $H_{23B}$ $C_{25B}$	120.1
C23A-	-C24A-C25A	120.20 (15)	$C_{23}B = C_{24}B = C_{25}B$	120.28 (15)
C23A-		119.9	C23B—C24B—H24B	119.9
C25A-		119.9	$C_{25B}$ $C_{24B}$ $H_{24B}$ $H_{24B}$	119.9
C20A-		119.85 (15)	$C_{24B} = C_{25B} = C_{20B}$	119.84 (14)
C20A-	-C25A-H25A	120.1	C24B—C25B—H25B	120.1
C24A-	-C25A-H25A	120.1	C20B—C25B—H25B	120.1
010A-		106.24 (15)	010B—C26B—C27B	108.21 (17)
010A-		110.15 (14)	O10B—C26B—C28B	107.62 (17)
C28A-		113.22 (17)	C27B—C26B—C28B	112.35 (16)
010A-		109.0	O10B—C26B—H26B	109.5
C28A-	C26AH26A	109.0	C27B—C26B—H26B	109.5
C27A-	C26AH26A	109.0	C28B—C26B—H26B	109.5
C26A-	—С27А—Н27А	109.5	C26B—C27B—H27D	109.5
C26A-		109.5	C26B—C27B—H27E	109.5
H27A-	—С27А—Н27В	109.5	H27D—C27B—H27E	109.5
C26A-	—С27А—Н27С	109.5	C26B—C27B—H27F	109.5
H27A-	—С27А—Н27С	109.5	H27D—C27B—H27F	109.5
H27B-	—С27А—Н27С	109.5	H27E—C27B—H27F	109.5
C26A-	C28AH28A	109.5	C26B—C28B—H28D	109.5

C26A—C28A—H28B	109.5	C26B—C28B—H28E	109.5
H28A—C28A—H28B	109.5	H28D—C28B—H28E	109.5
C26A—C28A—H28C	109.5	C26B—C28B—H28F	109.5
H28A—C28A—H28C	109.5	H28D—C28B—H28F	109.5
H28B—C28A—H28C	109.5	H28E—C28B—H28F	109.5
O11A—C29A—C31A	106.39 (13)	O11B—C29B—C31B	108.60 (14)
O11A-C29A-C30A	108.61 (14)	O11B-C29B-C30B	106.35 (17)
C31A—C29A—C30A	112.14 (17)	C31B—C29B—C30B	112.28 (17)
O11A—C29A—H29A	109.9	O11B—C29B—H29B	109.8
С31А—С29А—Н29А	109.9	C31B—C29B—H29B	109.8
C30A—C29A—H29A	109.9	C30B—C29B—H29B	109.8
C29A—C30A—H30A	109.5	C29B—C30B—H30D	109.5
C29A—C30A—H30B	109.5	C29B—C30B—H30E	109.5
H30A—C30A—H30B	109.5	H30D-C30B-H30E	109.5
C29A—C30A—H30C	109.5	C29B—C30B—H30F	109.5
H30A—C30A—H30C	109.5	H30D-C30B-H30F	109.5
H30B—C30A—H30C	109.5	H30E—C30B—H30F	109.5
C29A—C31A—H31A	109.5	C29B—C31B—H31D	109.5
C29A—C31A—H31B	109.5	C29B—C31B—H31E	109.5
H31A—C31A—H31B	109.5	H31D—C31B—H31E	109.5
C29A—C31A—H31C	109.5	C29B—C31B—H31F	109.5
H31A—C31A—H31C	109.5	H31D—C31B—H31F	109.5
H31B—C31A—H31C	109.5	H31E—C31B—H31F	109.5
O12A—C32A—C33A	108.69 (13)	O12B—C32B—C33B	109.34 (14)
O12A—C32A—C34A	106.76 (13)	O12B—C32B—C34B	106.40 (12)
C33A—C32A—C34A	113.43 (15)	C33B—C32B—C34B	112.50 (15)
O12A—C32A—H32A	109.3	O12B—C32B—H32B	109.5
C33A—C32A—H32A	109.3	C33B—C32B—H32B	109.5
C34A—C32A—H32A	109.3	C34B—C32B—H32B	109.5
С32А—С33А—Н33А	109.5	C32B—C33B—H33D	109.5
С32А—С33А—Н33В	109.5	C32B—C33B—H33E	109.5
H33A—C33A—H33B	109.5	H33D—C33B—H33E	109.5
C32A—C33A—H33C	109.5	C32B—C33B—H33F	109.5
H33A—C33A—H33C	109.5	H33D—C33B—H33F	109.5
H33B—C33A—H33C	109.5	H33E—C33B—H33F	109.5
C32A—C34A—H34A	109.5	C32B—C34B—H34D	109.5
C32A—C34A—H34B	109.5	C32B—C34B—H34E	109.5
H34A—C34A—H34B	109.5	H34D—C34B—H34E	109.5
C32A—C34A—H34C	109.5	C32B—C34B—H34F	109.5
H34A—C34A—H34C	109.5	H34D—C34B—H34F	109.5
H34B—C34A—H34C	109.5	H34E—C34B—H34F	109.5
O1A—C35A—Ru1A	171.58 (13)	O1B—C35B—Ru1B	174.05 (13)
O2A—C36A—Ru1A	177.84 (13)	O2B—C36B—Ru1B	179.52 (15)
O3A—C37A—Ru1A	172.82 (13)	O3B—C37B—Ru1B	173.51 (13)
O4A—C38A—Ru2A	175.30 (13)	O4B—C38B—Ru2B	173.16 (13)
O5A—C39A—Ru2A	176.23 (15)	O5B—C39B—Ru2B	176.98 (13)
O6A—C40A—Ru2A	172.77 (13)	O6B—C40B—Ru2B	173.57 (13)
O7A—C41A—Ru3A	173.30 (13)	O7B—C41B—Ru3B	174.79 (14)

O8A—C42A—Ru3A	178.35 (14)	O8B—C42B—Ru3B	175.89 (14)
O9A—C43A—Ru3A	172.86 (13)	O9B—C43B—Ru3B	174.57 (13)
	. ,		
C36A—Ru1A—Ru2A—C39A	67.2 (3)	C36B—Ru1B—Ru2B—C39B	-69.33 (19)
C37A—Ru1A—Ru2A—C39A	124.74 (14)	C35B—Ru1B—Ru2B—C39B	-138.77 (13)
C35A—Ru1A—Ru2A—C39A	-42.87 (14)	C37B—Ru1B—Ru2B—C39B	39.26 (13)
As1A—Ru1A—Ru2A—C39A	-141.84 (14)	As1B—Ru1B—Ru2B—C39B	130.64 (12)
Ru3A—Ru1A—Ru2A—C39A	31.40 (14)	Ru3B—Ru1B—Ru2B—C39B	-41.20(12)
C36A—Ru1A—Ru2A—C40A	134.0 (2)	C36B— $Ru1B$ — $Ru2B$ — $C38B$	-132.67(15)
C37A— $Ru1A$ — $Ru2A$ — $C40A$	-16841(6)	C35B = Ru1B = Ru2B = C38B	157 89 (6)
C35A— $Ru1A$ — $Ru2A$ — $C40A$	23.98 (6)	C37B $Ru1B$ $Ru2B$ $C38B$	-24.08(6)
$A_{s1}A$ $R_{u1}A$ $R_{u2}A$ $C40A$	-7499(4)	$A_{s1B}$ Ru1B Ru2B C38B	67 31 (5)
$R_{11}3A = R_{11}1A = R_{11}2A = C40A$	98 25 (4)	Ru3B Ru1B Ru2B C38B	-10454(5)
$C_{36} = R_{11} = R_{12} = C_{38}$	-45.2(2)	$C_{36}B_{\mu}B_{\mu}B_{\mu}B_{\mu}B_{\mu}C_{40}B_{\mu}B_{\mu}B_{\mu}B_{\mu}B_{\mu}B_{\mu}B_{\mu}B_{\mu$	44 09 (15)
$C_{37A} = Ru_{1A} = Ru_{2A} = C_{38A}$	+3.2(2) 12.38(6)	C35B $Bu1B$ $Bu2B$ $C40B$	-25.35(6)
$C_{37A}$ $Ru_{1A}$ $Ru_{2A}$ $C_{38A}$	-155 23 (6)	$C_{37B} = Ru_{1B} = Ru_{2B} = C_{40B}$	25.55 (0)
$A_{c1}A = B_{u1}A = B_{u2}A = C_{2}C_{2}A$	105.23(0)	$A_{c1}D = D_{u1}D = D_{u2}D = C40D$	-115.07(0)
ASIA— $Ru1A$ — $Ru2A$ — $C36ABu2A$ $Bu1A$ $Bu2A$ $C28A$	103.60 (4)	AS1D - Ku1B - Ku2D - C40B	-113.94(4)
Rusa - Rui A - Ruza - CsoA	-80.90(4)	Ru3D— $Ru1D$ — $Ru2D$ — $C40B$	12.22 (4)
$C_{30A}$ Rula Ruza Asza	-142.0(2)	C36B—RuIB—Ru2B—As2B	138.51(15)
$C_3/A$ —RuIA—Ru2A—As2A	-84.3/(4)	C35B—RuIB—Ru2B—As2B	69.07 (4)
C35A—RuIA—Ru2A—As2A	108.01 (4)	C3/B—RuIB—Ru2B—As2B	-112.91 (4)
AsIA—RuIA—Ru2A—As2A	9.043 (7)	As1B—Ru1B—Ru2B—As2B	-21.517 (6)
Ru3A—Ru1A—Ru2A—As2A	-17/.713 (6)	Ru3B—Ru1B—Ru2B—As2B	166.638 (6)
C36A—Ru1A—Ru2A—Ru3A	35.8 (2)	C36B—Ru1B—Ru2B—Ru3B	-28.13 (15)
C37A—Ru1A—Ru2A—Ru3A	93.34 (4)	C35B—Ru1B—Ru2B—Ru3B	-97.57 (4)
C35A—Ru1A—Ru2A—Ru3A	-74.27 (4)	C37B—Ru1B—Ru2B—Ru3B	80.46 (4)
As1A—Ru1A—Ru2A—Ru3A	-173.243 (6)	As1B—Ru1B—Ru2B—Ru3B	171.845 (6)
C39A—Ru2A—Ru3A—C42A	17.12 (7)	C39B—Ru2B—Ru3B—C42B	-23.60 (7)
C40A—Ru2A—Ru3A—C42A	109.72 (7)	C38B—Ru2B—Ru3B—C42B	-116.45 (6)
C38A—Ru2A—Ru3A—C42A	-75.28 (7)	C40B—Ru2B—Ru3B—C42B	64.79 (6)
As2A—Ru2A—Ru3A—C42A	-167.34 (5)	As2B—Ru2B—Ru3B—C42B	143.93 (5)
Ru1A—Ru2A—Ru3A—C42A	-173.44 (5)	Ru1B—Ru2B—Ru3B—C42B	170.65 (5)
C39A—Ru2A—Ru3A—C43A	109.02 (7)	C39B—Ru2B—Ru3B—C41B	-116.55 (7)
C40A—Ru2A—Ru3A—C43A	-158.38 (6)	C38B—Ru2B—Ru3B—C41B	150.60 (7)
C38A—Ru2A—Ru3A—C43A	16.62 (6)	C40B—Ru2B—Ru3B—C41B	-28.17 (7)
As2A—Ru2A—Ru3A—C43A	-75.44 (5)	As2B—Ru2B—Ru3B—C41B	50.97 (5)
Ru1A—Ru2A—Ru3A—C43A	-81.54 (5)	Ru1B—Ru2B—Ru3B—C41B	77.69 (5)
C39A—Ru2A—Ru3A—C41A	-72.97 (7)	C39B—Ru2B—Ru3B—C43B	70.57 (6)
C40A—Ru2A—Ru3A—C41A	19.63 (6)	C38B—Ru2B—Ru3B—C43B	-22.28(6)
C38A—Ru2A—Ru3A—C41A	-165.37 (6)	C40B—Ru2B—Ru3B—C43B	158.96 (6)
As2A—Ru2A—Ru3A—C41A	102.57 (5)	As2B—Ru2B—Ru3B—C43B	-121.90(5)
Ru1A—Ru2A—Ru3A—C41A	96.47 (4)	Ru1B—Ru2B—Ru3B—C43B	-95.18 (4)
C39A—Ru2A—Ru3A—P1A	-122.82 (7)	C39B—Ru2B—Ru3B—P1B	127.65 (5)
C40A—Ru2A—Ru3A—P1A	-30.21 (7)	C38B—Ru2B—Ru3B—P1B	34.80 (5)
C38A— $Ru2A$ — $Ru3A$ — $P1A$	144.78 (6)	C40B— $Ru2B$ — $Ru3B$ — $P1B$	-143.96(5)
As2A = Ru2A = Ru3A = P1A	52.72 (6)	$As_2B$ $Ru_2B$ $Ru_3B$ $P1B$	-64.83(3)
Ru1A Ru2A Ru3A P1A	46.62 (5)	Ru1B $Ru2B$ $Ru3B$ $P1B$	-38 11 (3)
C39A = Ru2A = Ru3A = Ru1A	-169.44(5)	C39B Ru2B Ru3B Ru3B Ru1B	165 76 (5)
Commuta Ruda Ruda	107. TT (3)	CJ/D_Ru2D_RuJD_Ru1D	105.70 (5)

C40A—Ru2A—Ru3A—Ru1A	-76.83 (4)	C38B—Ru2B—Ru3B—Ru1B	72.91 (4)
C38A—Ru2A—Ru3A—Ru1A	98.16 (4)	C40B—Ru2B—Ru3B—Ru1B	-105.85 (4)
As2A—Ru2A—Ru3A—Ru1A	6.104 (16)	As2B—Ru2B—Ru3B—Ru1B	-26.720 (11)
C36A—Ru1A—Ru3A—C42A	-160.05 (11)	C36B—Ru1B—Ru3B—C42B	149.67 (12)
C37A—Ru1A—Ru3A—C42A	-69.60 (11)	C35B—Ru1B—Ru3B—C42B	54.35 (12)
C35A—Ru1A—Ru3A—C42A	116.76 (11)	C37B—Ru1B—Ru3B—C42B	-123.23(12)
As1A—Ru1A—Ru3A—C42A	25.88 (10)	As1B—Ru1B—Ru3B—C42B	-39.62 (11)
Ru2A—Ru1A—Ru3A—C42A	12.94 (10)	Ru2B—Ru1B—Ru3B—C42B	-21.88 (11)
C36A—Ru1A—Ru3A—C43A	-72.77 (7)	C36B—Ru1B—Ru3B—C41B	69.08 (6)
C37A—Ru1A—Ru3A—C43A	17.68 (6)	C35B—Ru1B—Ru3B—C41B	-26.24 (6)
C35A—Ru1A—Ru3A—C43A	-155.96 (6)	C37B—Ru1B—Ru3B—C41B	156.18 (6)
As1A—Ru1A—Ru3A—C43A	113.15 (5)	As1B—Ru1B—Ru3B—C41B	-120.21 (5)
Ru2A—Ru1A—Ru3A—C43A	100.22 (4)	Ru2B—Ru1B—Ru3B—C41B	-102.47 (4)
C36A—Ru1A—Ru3A—C41A	104.41 (7)	C36B—Ru1B—Ru3B—C43B	-110.00 (6)
C37A—Ru1A—Ru3A—C41A	-165.14 (6)	C35B—Ru1B—Ru3B—C43B	154.68 (6)
C35A—Ru1A—Ru3A—C41A	21.22 (6)	C37B—Ru1B—Ru3B—C43B	-22.91 (6)
As1A—Ru1A—Ru3A—C41A	-69.67 (5)	As1B—Ru1B—Ru3B—C43B	60.71 (5)
Ru2A—Ru1A—Ru3A—C41A	-82.60 (5)	Ru2B—Ru1B—Ru3B—C43B	78.45 (4)
C36A—Ru1A—Ru3A—P1A	16.43 (5)	C36B—Ru1B—Ru3B—P1B	-22.22(5)
C37A—Ru1A—Ru3A—P1A	106.88 (4)	C35B—Ru1B—Ru3B—P1B	-117.54 (4)
C35A—Ru1A—Ru3A—P1A	-66.76 (4)	C37B—Ru1B—Ru3B—P1B	64.88 (4)
As1A—Ru1A—Ru3A—P1A	-157.642 (14)	As1B—Ru1B—Ru3B—P1B	148.491 (15)
Ru2A—Ru1A—Ru3A—P1A	-170.574 (12)	Ru2B—Ru1B—Ru3B—P1B	166.232 (10)
C36A—Ru1A—Ru3A—Ru2A	-172.99 (5)	C36B—Ru1B—Ru3B—Ru2B	171.55 (4)
C37A—Ru1A—Ru3A—Ru2A	-82.55 (4)	C35B—Ru1B—Ru3B—Ru2B	76.23 (4)
C35A—Ru1A—Ru3A—Ru2A	103.81 (4)	C37B—Ru1B—Ru3B—Ru2B	-101.35 (4)
As1A—Ru1A—Ru3A—Ru2A	12.933 (11)	As1B—Ru1B—Ru3B—Ru2B	-17.741 (12)
C36A—Ru1A—As1A—C7A	64.92 (7)	C36B—Ru1B—As1B—C7B	72.03 (6)
C37A—Ru1A—As1A—C7A	-26.33 (6)	C35B—Ru1B—As1B—C7B	167.47 (6)
C35A—Ru1A—As1A—C7A	153.85 (6)	C37B—Ru1B—As1B—C7B	-16.41 (6)
Ru2A—Ru1A—As1A—C7A	-109.66 (5)	Ru3B—Ru1B—As1B—C7B	-98.78 (5)
Ru3A—Ru1A—As1A—C7A	-120.61 (5)	Ru2B—Ru1B—As1B—C7B	-114.00 (5)
C36A—Ru1A—As1A—C1A	-50.90 (7)	C36B—Ru1B—As1B—C1B	-51.31 (7)
C37A—Ru1A—As1A—C1A	-142.14 (6)	C35B—Ru1B—As1B—C1B	44.12 (6)
C35A—Ru1A—As1A—C1A	38.04 (6)	C37B—Ru1B—As1B—C1B	-139.76 (6)
Ru2A—Ru1A—As1A—C1A	134.53 (5)	Ru3B—Ru1B—As1B—C1B	137.87 (5)
Ru3A—Ru1A—As1A—C1A	123.57 (5)	Ru2B—Ru1B—As1B—C1B	122.65 (5)
C36A—Ru1A—As1A—C13A	-172.07 (7)	C36B—Ru1B—As1B—C13B	-170.23 (7)
C37A—Ru1A—As1A—C13A	96.68 (6)	C35B—Ru1B—As1B—C13B	-74.79 (6)
C35A—Ru1A—As1A—C13A	-83.14 (6)	C37B—Ru1B—As1B—C13B	101.32 (6)
Ru2A—Ru1A—As1A—C13A	13.35 (5)	Ru3B—Ru1B—As1B—C13B	18.95 (5)
Ru3A—Ru1A—As1A—C13A	2.39 (5)	Ru2B—Ru1B—As1B—C13B	3.74 (5)
C39A—Ru2A—As2A—C14A	25.81 (7)	C39B—Ru2B—As2B—C14B	106.20 (7)
C40A—Ru2A—As2A—C14A	-65.50 (6)	C38B—Ru2B—As2B—C14B	-162.75 (7)
C38A—Ru2A—As2A—C14A	120.54 (6)	C40B—Ru2B—As2B—C14B	13.86 (7)
Ru3A—Ru2A—As2A—C14A	-149.73 (5)	Ru3B—Ru2B—As2B—C14B	-61.11 (6)
Ru1A—Ru2A—As2A—C14A	-144.33 (5)	Ru1B—Ru2B—As2B—C14B	-84.02 (5)
C39A—Ru2A—As2A—C20A	-99.44 (7)	C39B—Ru2B—As2B—C20B	-15.31 (7)

C40A—Ru2A—As2A—C20A	169.26 (6)	C38B—Ru2B—As2B—C20B	75.74 (6)
C38A—Ru2A—As2A—C20A	-4.70 (6)	C40B—Ru2B—As2B—C20B	-107.64 (6)
Ru3A—Ru2A—As2A—C20A	85.03 (5)	Ru3B—Ru2B—As2B—C20B	177.38 (5)
Ru1A—Ru2A—As2A—C20A	90.43 (5)	Ru1B—Ru2B—As2B—C20B	154.48 (5)
C39A—Ru2A—As2A—C13A	139.85 (7)	C39B—Ru2B—As2B—C13B	-131.51 (6)
C40A—Ru2A—As2A—C13A	48.54 (6)	C38B—Ru2B—As2B—C13B	-40.46 (6)
C38A—Ru2A—As2A—C13A	-125.42 (6)	C40B—Ru2B—As2B—C13B	136.16 (6)
Ru3A—Ru2A—As2A—C13A	-35.69 (5)	Ru3B—Ru2B—As2B—C13B	61.18 (5)
Ru1A—Ru2A—As2A—C13A	-30.29 (5)	Ru1B—Ru2B—As2B—C13B	38.28 (5)
C42A—Ru3A—P1A—O11A	-116.15 (7)	C42B—Ru3B—P1B—O12B	127.58 (7)
C43A—Ru3A—P1A—O11A	151.78 (6)	C41B—Ru3B—P1B—O12B	-139.36(7)
C41A—Ru3A—P1A—O11A	-26.61 (6)	C43B—Ru3B—P1B—O12B	32.98 (7)
Ru2A—Ru3A—P1A—O11A	23.09 (8)	Ru2B—Ru3B—P1B—O12B	-22.73 (6)
Ru1A—Ru3A—P1A—O11A	65.67 (5)	Ru1B—Ru3B—P1B—O12B	-56.05 (5)
C42A—Ru3A—P1A—O10A	-2.23 (7)	C42B—Ru3B—P1B—O11B	-116.72 (7)
C43A—Ru3A—P1A—O10A	-94.30 (7)	C41B—Ru3B—P1B—O11B	-23.67 (7)
C41A—Ru3A—P1A—O10A	87.30 (7)	C43B—Ru3B—P1B—O11B	148.68 (7)
Ru2A—Ru3A—P1A—O10A	137.01 (6)	Ru2B—Ru3B—P1B—O11B	92.97 (6)
Ru1A—Ru3A—P1A—O10A	179.58 (5)	Ru1B—Ru3B—P1B—O11B	59.65 (5)
C42A—Ru3A—P1A—O12A	117.14 (7)	C42B—Ru3B—P1B—O10B	5.80 (8)
C43A—Ru3A—P1A—O12A	25.07 (7)	C41B—Ru3B—P1B—O10B	98.85 (7)
C41A—Ru3A—P1A—O12A	-153.32 (7)	C43B—Ru3B—P1B—O10B	-88.80 (7)
Ru2A—Ru3A—P1A—O12A	-103.62 (7)	Ru2B—Ru3B—P1B—O10B	-144.51 (6)
Ru1A—Ru3A—P1A—O12A	-61.04 (5)	Ru1B—Ru3B—P1B—O10B	-177.83 (6)
O11A—P1A—O10A—C26A	179.62 (11)	O12B—P1B—O10B—C26B	-89.98 (14)
O12A—P1A—O10A—C26A	-70.77 (12)	O11B—P1B—O10B—C26B	166.78 (13)
Ru3A—P1A—O10A—C26A	60.09 (12)	Ru3B—P1B—O10B—C26B	33.39 (15)
O10A—P1A—O11A—C29A	33.16 (13)	O12B—P1B—O11B—C29B	-166.45 (12)
O12A—P1A—O11A—C29A	-68.72 (12)	O10B—P1B—O11B—C29B	-58.39 (13)
Ru3A—P1A—O11A—C29A	157.56 (10)	Ru3B—P1B—O11B—C29B	73.39 (13)
O11A—P1A—O12A—C32A	-79.79 (12)	O11B—P1B—O12B—C32B	50.58 (13)
O10A—P1A—O12A—C32A	176.79 (12)	O10B—P1B—O12B—C32B	-51.39 (13)
Ru3A—P1A—O12A—C32A	48.22 (13)	Ru3B—P1B—O12B—C32B	179.28 (11)
C7A—As1A—C1A—C6A	83.82 (13)	C7B—As1B—C1B—C6B	-132.60 (12)
C13A—As1A—C1A—C6A	-20.86 (14)	C13B—As1B—C1B—C6B	122.97 (12)
Ru1A—As1A—C1A—C6A	-149.85 (11)	Ru1B—As1B—C1B—C6B	-3.14 (13)
C7A—As1A—C1A—C2A	-89.09 (12)	C7B—As1B—C1B—C2B	52.11 (13)
C13A—As1A—C1A—C2A	166.23 (12)	C13B—As1B—C1B—C2B	-52.32 (13)
Ru1A—As1A—C1A—C2A	37.25 (13)	Ru1B—As1B—C1B—C2B	-178.43 (10)
C6A—C1A—C2A—C3A	-0.7 (2)	C6B—C1B—C2B—C3B	-1.5 (2)
As1A—C1A—C2A—C3A	172.34 (13)	As1B—C1B—C2B—C3B	173.85 (12)
C1A—C2A—C3A—C4A	0.9 (3)	C1B—C2B—C3B—C4B	0.0 (2)
C2A—C3A—C4A—C5A	0.2 (3)	C2B—C3B—C4B—C5B	1.5 (2)
C3A—C4A—C5A—C6A	-1.4 (3)	C3B—C4B—C5B—C6B	-1.4 (2)
C2A-C1A-C6A-C5A	-0.5 (2)	C2B—C1B—C6B—C5B	1.6 (2)
As1A—C1A—C6A—C5A	-173.35 (13)	As1B—C1B—C6B—C5B	-173.71 (11)
C4A—C5A—C6A—C1A	1.6 (3)	C4B—C5B—C6B—C1B	-0.1 (2)
C1A—As1A—C7A—C8A	-75.93 (13)	C1B—As1B—C7B—C8B	-145.53 (14)

C13A—As1A—C7A—C8A	25.92 (14)	C13B—As1B—C7B—C8B	-42.27 (15)
Ru1A—As1A—C7A—C8A	155.49 (11)	Ru1B—As1B—C7B—C8B	82.55 (14)
C1A—As1A—C7A—C12A	98.57 (12)	C1B—As1B—C7B—C12B	42.33 (13)
C13A—As1A—C7A—C12A	-159.59 (11)	C13B—As1B—C7B—C12B	145.60 (12)
Ru1A—As1A—C7A—C12A	-30.01 (13)	Ru1B—As1B—C7B—C12B	-89.59 (12)
C12A—C7A—C8A—C9A	0.3 (2)	C12B—C7B—C8B—C9B	-1.0(3)
As1A—C7A—C8A—C9A	174.73 (12)	As1B—C7B—C8B—C9B	-173.15 (15)
C7A—C8A—C9A—C10A	-0.1 (3)	C7B-C8B-C9B-C10B	0.3 (3)
C8A—C9A—C10A—C11A	-0.2 (3)	C8B—C9B—C10B—C11B	0.6 (3)
C9A—C10A—C11A—C12A	0.3 (2)	C9B—C10B—C11B—C12B	-0.9(3)
C10A—C11A—C12A—C7A	-0.1 (2)	C10B—C11B—C12B—C7B	0.1 (2)
C8A—C7A—C12A—C11A	-0.2 (2)	C8B—C7B—C12B—C11B	0.8 (2)
As1A—C7A—C12A—C11A	-174.87 (12)	As1B—C7B—C12B—C11B	173.03 (12)
C14A—As2A—C13A—As1A	165.10 (7)	C14B—As2B—C13B—As1B	88.38 (8)
C20A—As2A—C13A—As1A	-88.44 (8)	C20B—As2B—C13B—As1B	-168.92(7)
Ru2A—As2A—C13A—As1A	43.52 (7)	Ru2B—As2B—C13B—As1B	-42.52 (8)
C7A—As1A—C13A—As2A	93.09 (7)	C7B—As1B—C13B—As2B	147.51 (7)
C1A—As1A—C13A—As2A	-168.20(7)	C1B—As1B—C13B—As2B	-107.27 (8)
Ru1A—As1A—C13A—As2A	-36.83 (8)	Ru1B—As1B—C13B—As2B	22.10 (9)
C20A—As2A—C14A—C19A	-108.23 (13)	C20B—As2B—C14B—C19B	-107.22 (13)
C13A—As2A—C14A—C19A	-2.02 (13)	C13B—As2B—C14B—C19B	-3.33 (14)
Ru2A—As2A—C14A—C19A	116.80 (11)	Ru2B—As2B—C14B—C19B	122.33 (12)
C20A—As2A—C14A—C15A	75.49 (12)	C20B—As2B—C14B—C15B	68.55 (12)
C13A—As2A—C14A—C15A	-178.31 (12)	C13B—As2B—C14B—C15B	172.44 (11)
Ru2A—As2A—C14A—C15A	-59.49 (12)	Ru2B—As2B—C14B—C15B	-61.90 (12)
C19A—C14A—C15A—C16A	1.1 (2)	C19B—C14B—C15B—C16B	1.5 (2)
As2A—C14A—C15A—C16A	177.56 (13)	As2B—C14B—C15B—C16B	-174.46 (13)
C14A—C15A—C16A—C17A	-0.7 (3)	C14B—C15B—C16B—C17B	-1.0 (3)
C15A—C16A—C17A—C18A	-0.3 (3)	C15B—C16B—C17B—C18B	-0.5(3)
C16A—C17A—C18A—C19A	0.8 (2)	C16B—C17B—C18B—C19B	1.5 (3)
C15A—C14A—C19A—C18A	-0.6 (2)	C15B—C14B—C19B—C18B	-0.5 (2)
As2A—C14A—C19A—C18A	-176.79 (11)	As2B-C14B-C19B-C18B	175.18 (13)
C17A—C18A—C19A—C14A	-0.4 (2)	C17B—C18B—C19B—C14B	-1.0(3)
C14A—As2A—C20A—C21A	-174.38 (11)	C14B—As2B—C20B—C21B	18.73 (14)
C13A—As2A—C20A—C21A	79.94 (12)	C13B—As2B—C20B—C21B	-88.01 (14)
Ru2A—As2A—C20A—C21A	-44.27 (13)	Ru2B—As2B—C20B—C21B	150.28 (12)
C14A—As2A—C20A—C25A	5.91 (13)	C14B—As2B—C20B—C25B	-159.07 (12)
C13A—As2A—C20A—C25A	-99.77 (12)	C13B—As2B—C20B—C25B	94.18 (13)
Ru2A—As2A—C20A—C25A	136.02 (10)	Ru2B—As2B—C20B—C25B	-27.52 (14)
C25A—C20A—C21A—C22A	1.2 (2)	C25B—C20B—C21B—C22B	0.5 (3)
As2A—C20A—C21A—C22A	-178.53 (12)	As2B-C20B-C21B-C22B	-177.27 (14)
C20A—C21A—C22A—C23A	-1.3 (2)	C20B—C21B—C22B—C23B	-0.3 (3)
C21A—C22A—C23A—C24A	0.3 (3)	C21B—C22B—C23B—C24B	-0.4 (3)
C22A—C23A—C24A—C25A	0.8 (3)	C22B—C23B—C24B—C25B	0.9 (3)
C21A—C20A—C25A—C24A	-0.1 (2)	C23B—C24B—C25B—C20B	-0.7 (3)
As2A—C20A—C25A—C24A	179.66 (12)	C21B—C20B—C25B—C24B	0.0 (2)
C23A—C24A—C25A—C20A	-1.0 (2)	As2B-C20B-C25B-C24B	177.84 (13)
P1A-010A-C26A-C28A	-147.83 (13)	P1B-010B-C26B-C27B	116.73 (17)

P1A-010A-C26A-C27A	89.19 (18)	P1B-010B-C26B-C28B	-121.64 (15)
P1A-011A-C29A-C31A	143.59 (14)	P1B-011B-C29B-C31B	101.12 (17)
P1A-011A-C29A-C30A	-95.51 (17)	P1B-011B-C29B-C30B	-137.85 (13)
P1A-012A-C32A-C33A	90.92 (14)	P1B-012B-C32B-C33B	-98.61 (15)
P1A—O12A—C32A—C34A	-146.38(12)	P1B-012B-C32B-C34B	139.65 (13)
C36A—Ru1A—C35A—O1A	7.7 (9)	C36B—Ru1B—C35B—O1B	48.2 (12)
C37A—Ru1A—C35A—O1A	89.9 (10)	C37B—Ru1B—C35B—O1B	-174.1 (10)
As1A—Ru1A—C35A—O1A	-90.9 (9)	As1B—Ru1B—C35B—O1B	-54.2 (12)
Ru2A—Ru1A—C35A—O1A	177 (100)	Ru3B—Ru1B—C35B—O1B	153.4 (12)
Ru3A—Ru1A—C35A—O1A	121.2 (9)	Ru2B—Ru1B—C35B—O1B	-148.2(12)
C37A—Ru1A—C36A—O2A	70 (4)	C35B—Ru1B—C36B—O2B	-100 (15)
C35A—Ru1A—C36A—O2A	-122 (4)	C37B—Ru1B—C36B—O2B	83 (15)
As1A—Ru1A—C36A—O2A	-24 (4)	As1B—Ru1B—C36B—O2B	-8 (15)
Ru2A—Ru1A—C36A—O2A	127 (4)	Ru3B—Ru1B—C36B—O2B	168 (100)
Ru3A—Ru1A—C36A—O2A	160 (4)	Ru2B—Ru1B—C36B—O2B	-167(100)
C36A—Ru1A—C37A—O3A	-21.1 (11)	C36B—Ru1B—C37B—O3B	-42.7 (12)
C35A—Ru1A—C37A—O3A	-103.0(11)	C35B—Ru1B—C37B—O3B	179 (100)
As1A—Ru1A—C37A—O3A	77.8 (11)	As1B—Ru1B—C37B—O3B	59.5 (12)
Ru2A—Ru1A—C37A—O3A	168.2 (11)	Ru3B—Ru1B—C37B—O3B	-148.0(12)
Ru3A—Ru1A—C37A—O3A	-133.8(11)	Ru2B—Ru1B—C37B—O3B	153.9 (12)
C39A—Ru2A—C38A—O4A	42.3 (16)	C39B—Ru2B—C38B—O4B	49.3 (12)
C40A—Ru2A—C38A—O4A	-163.5 (14)	C40B—Ru2B—C38B—O4B	167.8 (9)
As2A—Ru2A—C38A—O4A	-59.1 (16)	As2B—Ru2B—C38B—O4B	-57.5 (12)
Ru3A—Ru2A—C38A—O4A	142.9 (16)	Ru3B—Ru2B—C38B—O4B	152.9 (11)
Ru1A—Ru2A—C38A—O4A	-156.3 (16)	Ru1B—Ru2B—C38B—O4B	-149.6 (12)
C40A—Ru2A—C39A—O5A	-6 (2)	C38B—Ru2B—C39B—O5B	34 (3)
C38A—Ru2A—C39A—O5A	172 (2)	C40B—Ru2B—C39B—O5B	-142(3)
As2A—Ru2A—C39A—O5A	-91 (2)	As2B—Ru2B—C39B—O5B	124 (3)
Ru3A—Ru2A—C39A—O5A	87 (2)	Ru3B—Ru2B—C39B—O5B	-63 (3)
Ru1A—Ru2A—C39A—O5A	59 (2)	Ru1B—Ru2B—C39B—O5B	-27 (3)
C39A—Ru2A—C40A—O6A	-62.4 (11)	C39B—Ru2B—C40B—O6B	-35.6 (12)
C38A—Ru2A—C40A—O6A	143.4 (9)	C38B—Ru2B—C40B—O6B	-154.0 (10)
As2A—Ru2A—C40A—O6A	38.4 (11)	As2B—Ru2B—C40B—O6B	71.4 (12)
Ru3A—Ru2A—C40A—O6A	-163.4 (11)	Ru3B—Ru2B—C40B—O6B	-139.0 (12)
Ru1A—Ru2A—C40A—O6A	136.1 (11)	Ru1B—Ru2B—C40B—O6B	163.9 (12)
C42A—Ru3A—C41A—O7A	46.8 (12)	C42B—Ru3B—C41B—O7B	44.8 (16)
C43A—Ru3A—C41A—O7A	-86.7 (18)	C43B—Ru3B—C41B—O7B	-153.8 (14)
P1A—Ru3A—C41A—O7A	-52.9 (12)	P1B—Ru3B—C41B—O7B	-59.9 (16)
Ru2A—Ru3A—C41A—O7A	136.4 (12)	Ru2B—Ru3B—C41B—O7B	140.0 (16)
Ru1A—Ru3A—C41A—O7A	-163.7 (12)	Ru1B—Ru3B—C41B—O7B	-160.7 (16)
C43A—Ru3A—C42A—O8A	-106 (6)	C41B—Ru3B—C42B—O8B	59 (2)
C41A—Ru3A—C42A—O8A	76 (6)	C43B—Ru3B—C42B—O8B	-119 (2)
P1A—Ru3A—C42A—O8A	163 (6)	P1B—Ru3B—C42B—O8B	152 (2)
Ru2A—Ru3A—C42A—O8A	-9 (6)	Ru2B—Ru3B—C42B—O8B	-39 (2)
Ru1A—Ru3A—C42A—O8A	-21 (6)	Ru1B—Ru3B—C42B—O8B	-20 (2)
C42A—Ru3A—C43A—O9A	-57.8 (12)	C42B—Ru3B—C43B—O9B	-66.2 (15)
C41A—Ru3A—C43A—O9A	75.7 (19)	C41B—Ru3B—C43B—O9B	132.3 (13)
P1A—Ru3A—C43A—O9A	42.0 (12)	P1B—Ru3B—C43B—O9B	38.2 (15)

Ru2A—Ru3A—C43A—O9A	-147.6 (12)	Ru2B—Ru3B—C43B—O9B	-160.3 (15)
Ru1A—Ru3A—C43A—O9A	152.9 (12)	Ru1B—Ru3B—C43B—O9B	139.2 (15)

Hydrogen-bond geometry (Å, °)

D—H···A	D—H	H…A	D····A	<i>D</i> —H··· <i>A</i>
C10 <i>B</i> —H10 <i>B</i> ····O6 <i>A</i> <sup>i</sup>	0.93	2.59	3.293 (2)	132
C17 <i>B</i> —H17 <i>B</i> ····O2 <i>B</i> <sup>ii</sup>	0.93	2.57	3.477 (2)	165
$C33A$ — $H33B$ ···O2 $A^{iii}$	0.96	2.56	3.492 (2)	164

Symmetry codes: (i) -*x*, -*y*+1, -*z*; (ii) -*x*+1, -*y*+1, -*z*; (iii) -*x*, -*y*+1, -*z*+1.