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(Dithiobenzoato- κ^2S,S')[hydridotris(pyrazol-1-yl- κN^2)borato](triphenylphosphine- κP)ruthenium(II)

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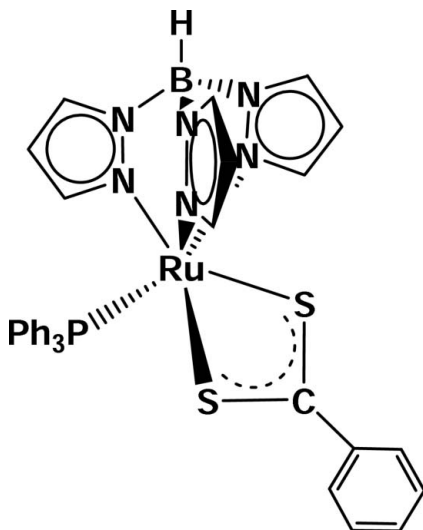
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 Key indicators: single-crystal X-ray study; $T = 200$ K; mean $\sigma(C-C) = 0.005$ Å; R factor = 0.036; wR factor = 0.080; data-to-parameter ratio = 13.7.

Reaction of $[Ru(Tp)Cl(PPh_3)_2]$ ($Tp =$ hydridotrispyrazolylborate) with ammonium dithiobenzoate in methanol leads to the formation of the title compound, $[Ru(C_9H_{10}BN_6)(C_7H_5S_2)(C_{18}H_{15}P)]$. In the crystal structure, the Ru atom is coordinated by three N atoms of the Tp ligand, one P atom of the triphenylphosphine ligand and the two S atoms of the dithiobenzoate ligand within a slightly distorted octahedron. The Ru–S bonds are slightly different [2.321 (1) and 2.396 (1) Å] and the average N–Ru–N angle is 86.31°.

Related literature

For general background, see: Alock *et al.* (1992); Burrows (2001); Pavlik *et al.* (2005); Hidai *et al.* (2000); Vit & Zdrzil (1989). For related structures, see: Gemel *et al.* (1996); Slugovc *et al.* (1998); Sellmann *et al.* (1999); Meno *et al.* (1995).



Experimental

Crystal data

$[Ru(C_9H_{10}BN_6)(C_7H_5S_2)(C_{18}H_{15}P)]$	$V = 3184.2$ (6) Å ³
$M_r = 729.61$	$Z = 4$
Monoclinic, $P2_1/n$	Mo $K\alpha$ radiation
$a = 12.8915$ (13) Å	$\mu = 0.71$ mm ⁻¹
$b = 18.394$ (2) Å	$T = 200$ (2) K
$c = 13.5174$ (16) Å	$0.18 \times 0.12 \times 0.02$ mm
$\beta = 96.591$ (5)°	

Data collection

Nonius KappaCCD diffractometer	22786 measured reflections
Absorption correction: multi-scan (Blessing, 1995)	5557 independent reflections
$T_{min} = 0.883$, $T_{max} = 0.986$	3974 reflections with $I > 2\sigma(I)$
	$R_{int} = 0.056$

Refinement

$R[F^2 > 2\sigma(F^2)] = 0.036$	406 parameters
$wR(F^2) = 0.080$	H-atom parameters constrained
$S = 1.02$	$\Delta\rho_{max} = 1.39$ e Å ⁻³
5557 reflections	$\Delta\rho_{min} = -0.62$ e Å ⁻³

Data collection: *COLLECT* (Nonius, 1999); cell refinement: *DENZO* and *SCALEPACK* (Otwinowski & Minor, 1997); data reduction: *DENZO* and *SCALEPACK*; program(s) used to solve structure: *SHELXS97* (Sheldrick, 2008); program(s) used to refine structure: *SHELXL97* (Sheldrick, 2008); molecular graphics: *ORTEP-3 for Windows* (Farrugia, 1997); software used to prepare material for publication: *WinGX* (Farrugia, 1999).

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Supplementary data and figures for this paper are available from the IUCr electronic archives (Reference: NC2121).

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supporting information

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(Dithiobenzoato- κ^2S,S')[hydridotris(pyrazol-1-yl- κN^2)borato](triphenylphosphine- κP)ruthenium(II)

Chia-Her Lin, Yao-Ren Liang, Hung-Chun Tong, Yih Hsing Lo and Ting Shen Kuo

S1. Comment

Ruthenium(II) hydridotripyrazolylborate complexes, (Ru(Tp)), are of interest for stoichiometric and catalytic transformations of organic molecules (Pavlik *et al.*, 2005). The complex [Ru(Tp)Cl(PPh₃)₂] (Alock *et al.*, 1992) has been used as the starting material for the synthesis of several complexes because the chloride atom and the phosphine ligand can easily be substituted (Burrows, 2001). On the other hand, the chemistry of transition metal sulfur compounds has attracted much interest for their importance in the field of catalysis and metalloenzymes (Hidai *et al.*, 2000). In recent years there has been an increased interest in ruthenium sulfur complexes, in part because of the high catalytic activity of RuS₂ in various hydrotreating processes (Vit & Zdrzil, 1989). Thus, many ruthenium thiolate complexes have been reported. However, ruthenium complexes with dithio ligands are relatively rare (Sellmann *et al.*, 1999). In this context the structure of the title compound was determined.

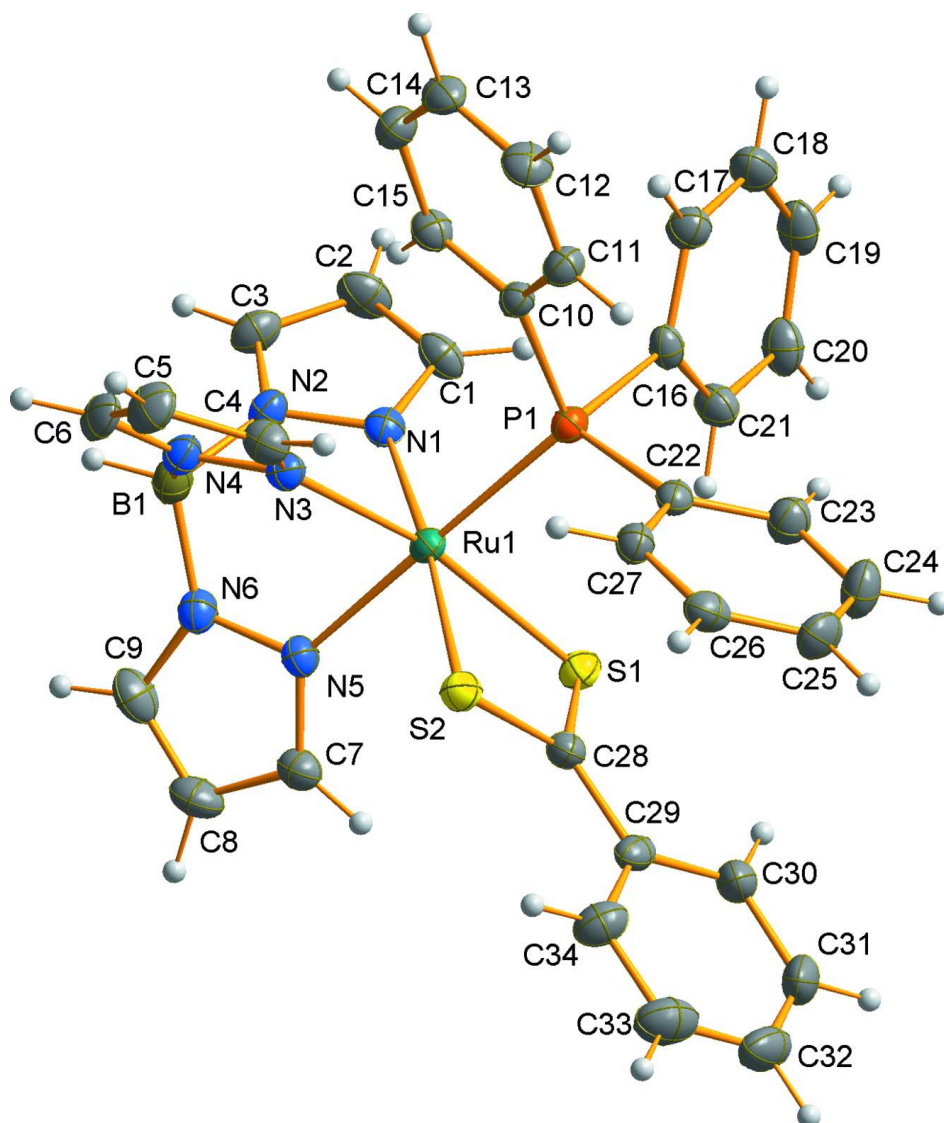
In the crystal structure of the title compound, the Ru atom is coordinated by three N atoms of the Tp ligand, two S atoms of the dithiobenzoate ligand and one P atom of the triphenylphosphine ligand forming slightly distorted octahedron. The average N—Ru—N angle amount to 86.31° and the three Ru—N bond lengths of 2.141 (3), 2.098 (3) and 2.134 (3) Å are slightly longer than the average distance of 2.038 Å in observed in other RuTp complexes (Gemel *et al.* 1996 and Slugovc *et al.* 1998). The dithiobenzoate ligand chelates the ruthenium centre with slightly different Ru—S bonds of 2.321 (1) Å and 2.396 (1) Å and an S—Ru—S angle of 71.61 (3)°. The average Ru—S bond length of 2.3588 (11) Å is slightly shorter than in *cis*-[Ru(S₂CNET₂)₂(PPh₃)₂] (av. 2.3952 (5) Å) (Meno *et al.*, 1995).

S2. Experimental

To a solution of [Ru(Tp)Cl(PPh₃)₂] (3.95 g, 4.50 mmol) in MeOH (20 ml) an excess of [NH₄][S₂C(C₆H₅)] (1.71 g, 10 mmol) were added. The reaction mixture was stirred for 4 h at room temperature. The solvent was removed in vacuum and 20 ml of CH₂Cl₂ were added to the residue. After filtration the solvent was removed in vacuum to give the title compound. Spectroscopic analysis: IR(KBr, cm⁻¹): ν (BH)2467 cm⁻¹. ³¹P NMR(CDCl₃, 303 K, δ ,p.p.m.): d 58.3 (PPh₃). MS (*m/z*,Ru102): 730.2 (*M*⁺), 468.1 (*M*⁺ - PPh₃). Anal. Calcd for C₃₄H₃₀BN₆PRuS₂: C, 55.97; H,4.14; N, 11.52. Found: C, 55.73; H, 4.11;N, 11.42. The bright-yellow crystals used for X-ray structure analysis were obtained by recrystallization of the crude product from dichloromethane–hexane.

S3. Refinement

The H atoms were placed in idealized positions and constrained to ride on their parent atoms, with C—H = 0.95 Å and $U_{iso}(H) = 1.2U_{eq}(C)$, B—H = 1.0 Å and $U_{iso}(H) = 1.2U_{eq}(B)$.

**Figure 1**

Molecular structure of the title compound with labelling and displacement ellipsoids drawn at the 30% probability level (H atoms are shown as spheres of arbitrary radius).

(Dithiobenzoato- κ^2S,S')[hydridotris(pyrazol-1-yl- κN^2)borato](triphenylphosphine- κP)ruthenium(II)

Crystal data

[Ru(C₉H₁₀BN₆)(C₇H₅S₂)(C₁₈H₁₅P)]

$M_r = 729.61$

Monoclinic, $P2_1/n$

Hall symbol: -P 2yn

$a = 12.8915 (13) \text{ \AA}$

$b = 18.394 (2) \text{ \AA}$

$c = 13.5174 (16) \text{ \AA}$

$\beta = 96.591 (5)^\circ$

$V = 3184.2 (6) \text{ \AA}^3$

$Z = 4$

$F(000) = 1488$

$D_x = 1.522 \text{ Mg m}^{-3}$

Mo $K\alpha$ radiation, $\lambda = 0.71073 \text{ \AA}$

Cell parameters from 5326 reflections

$\theta = 2.3\text{--}24.9^\circ$

$\mu = 0.71 \text{ mm}^{-1}$

$T = 200 \text{ K}$

Prism, green

$0.18 \times 0.12 \times 0.02 \text{ mm}$

Data collection

Nonius KappaCCD diffractometer	22786 measured reflections
Radiation source: fine-focus sealed tube	5557 independent reflections
Graphite monochromator	3974 reflections with $I > 2\sigma(I)$
CCD rotation images, thick slices scans	$R_{\text{int}} = 0.056$
Absorption correction: multi-scan (Blessing, 1995)	$\theta_{\text{max}} = 25.0^\circ$, $\theta_{\text{min}} = 1.9^\circ$
$T_{\text{min}} = 0.883$, $T_{\text{max}} = 0.986$	$h = -15 \rightarrow 12$
	$k = -21 \rightarrow 20$
	$l = -16 \rightarrow 15$

Refinement

Refinement on F^2	Secondary atom site location: difference Fourier map
Least-squares matrix: full	Hydrogen site location: inferred from neighbouring sites
$R[F^2 > 2\sigma(F^2)] = 0.036$	H-atom parameters constrained
$wR(F^2) = 0.080$	$w = 1/[\sigma^2(F_o^2) + (0.0315P)^2 + 1.684P]$
$S = 1.02$	where $P = (F_o^2 + 2F_c^2)/3$
5557 reflections	$(\Delta/\sigma)_{\text{max}} = 0.001$
406 parameters	$\Delta\rho_{\text{max}} = 1.39 \text{ e } \text{\AA}^{-3}$
0 restraints	$\Delta\rho_{\text{min}} = -0.62 \text{ e } \text{\AA}^{-3}$
Primary atom site location: structure-invariant direct methods	

Special details

Experimental. Semi-empirical from equivalents by WinGX (Blessing, 1995)

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

	<i>x</i>	<i>y</i>	<i>z</i>	$U_{\text{iso}}^*/U_{\text{eq}}$
B1	0.3330 (3)	0.7338 (2)	0.0521 (3)	0.0277 (11)
H1'	0.3875	0.7068	0.0198	0.033*
C1	0.3375 (3)	0.75212 (19)	0.3191 (3)	0.0272 (9)
H1	0.3096	0.7678	0.3777	0.033*
C2	0.4285 (3)	0.7130 (2)	0.3191 (3)	0.0331 (10)
H2	0.4739	0.6970	0.3752	0.040*
C3	0.4393 (3)	0.7023 (2)	0.2210 (3)	0.0283 (9)
H3	0.4949	0.6770	0.1959	0.034*
C4	0.2720 (3)	0.9240 (2)	0.0120 (3)	0.0282 (9)
H4	0.2359	0.9673	0.0255	0.034*
C5	0.3432 (3)	0.9186 (2)	-0.0566 (3)	0.0334 (10)
H5	0.3648	0.9559	-0.0983	0.040*
C6	0.3760 (3)	0.8478 (2)	-0.0514 (3)	0.0310 (10)
H6	0.4256	0.8267	-0.0897	0.037*
C7	0.0588 (3)	0.68809 (19)	0.0294 (3)	0.0259 (9)

H7	-0.0109	0.6970	0.0424	0.031*
C8	0.0901 (3)	0.6302 (2)	-0.0258 (3)	0.0338 (10)
H8	0.0474	0.5927	-0.0570	0.041*
C9	0.1945 (3)	0.6385 (2)	-0.0257 (3)	0.0327 (10)
H9	0.2389	0.6070	-0.0574	0.039*
C10	0.2846 (3)	0.98437 (19)	0.2445 (2)	0.0191 (8)
C11	0.2725 (3)	1.05912 (19)	0.2465 (3)	0.0243 (9)
H11	0.2057	1.0794	0.2525	0.029*
C12	0.3569 (3)	1.1045 (2)	0.2400 (3)	0.0298 (10)
H12	0.3478	1.1557	0.2408	0.036*
C13	0.4542 (3)	1.0757 (2)	0.2322 (3)	0.0279 (9)
H13	0.5124	1.1070	0.2287	0.034*
C14	0.4670 (3)	1.0014 (2)	0.2297 (3)	0.0280 (9)
H14	0.5341	0.9814	0.2242	0.034*
C15	0.3825 (3)	0.9555 (2)	0.2350 (3)	0.0241 (9)
H15	0.3916	0.9043	0.2323	0.029*
C16	0.2187 (3)	0.89959 (19)	0.3980 (3)	0.0218 (8)
C17	0.2997 (3)	0.9362 (2)	0.4538 (3)	0.0275 (9)
H17	0.3328	0.9756	0.4247	0.033*
C18	0.3329 (3)	0.9164 (2)	0.5511 (3)	0.0349 (10)
H18	0.3890	0.9418	0.5877	0.042*
C19	0.2853 (3)	0.8606 (2)	0.5942 (3)	0.0329 (10)
H19	0.3102	0.8460	0.6601	0.040*
C20	0.2015 (3)	0.8251 (2)	0.5433 (3)	0.0322 (10)
H20	0.1661	0.7878	0.5747	0.039*
C21	0.1693 (3)	0.8448 (2)	0.4452 (3)	0.0273 (9)
H21	0.1120	0.8200	0.4096	0.033*
C22	0.0647 (3)	0.97588 (18)	0.2688 (3)	0.0206 (8)
C23	0.0053 (3)	0.9746 (2)	0.3480 (3)	0.0301 (10)
H23	0.0297	0.9485	0.4068	0.036*
C24	-0.0893 (3)	1.0110 (2)	0.3421 (3)	0.0380 (11)
H24	-0.1290	1.0100	0.3971	0.046*
C25	-0.1262 (3)	1.0486 (2)	0.2574 (3)	0.0341 (10)
H25	-0.1921	1.0722	0.2530	0.041*
C26	-0.0671 (3)	1.0517 (2)	0.1791 (3)	0.0279 (9)
H26	-0.0913	1.0787	0.1212	0.034*
C27	0.0272 (3)	1.01580 (19)	0.1842 (3)	0.0228 (9)
H27	0.0673	1.0182	0.1295	0.027*
C28	-0.0643 (3)	0.83083 (19)	0.1494 (3)	0.0212 (8)
C29	-0.1777 (3)	0.83849 (19)	0.1434 (3)	0.0256 (9)
C30	-0.2313 (3)	0.8185 (2)	0.2228 (3)	0.0341 (10)
H30	-0.1940	0.8001	0.2822	0.041*
C31	-0.3392 (3)	0.8253 (3)	0.2156 (3)	0.0454 (12)
H31	-0.3755	0.8108	0.2697	0.055*
C32	-0.3936 (3)	0.8529 (2)	0.1303 (4)	0.0457 (12)
H32	-0.4673	0.8575	0.1258	0.055*
C33	-0.3420 (3)	0.8735 (2)	0.0523 (4)	0.0452 (12)
H33	-0.3801	0.8927	-0.0063	0.054*

C34	-0.2346 (3)	0.8668 (2)	0.0578 (3)	0.0334 (10)
H34	-0.1993	0.8815	0.0032	0.040*
N1	0.2933 (2)	0.76531 (15)	0.2267 (2)	0.0215 (7)
N2	0.3582 (2)	0.73352 (16)	0.1657 (2)	0.0235 (7)
N3	0.2606 (2)	0.86080 (16)	0.0565 (2)	0.0210 (7)
N4	0.3261 (2)	0.81318 (16)	0.0171 (2)	0.0241 (7)
N5	0.1408 (2)	0.72945 (15)	0.0615 (2)	0.0223 (7)
N6	0.2250 (2)	0.69857 (16)	0.0266 (2)	0.0251 (7)
P1	0.18113 (7)	0.91954 (5)	0.26504 (7)	0.0190 (2)
Ru1	0.16002 (2)	0.822395 (15)	0.15631 (2)	0.01805 (9)
S1	0.01327 (7)	0.78411 (5)	0.23709 (7)	0.0227 (2)
S2	0.00968 (7)	0.87056 (5)	0.06980 (7)	0.0231 (2)

Atomic displacement parameters (Å²)

	U^{11}	U^{22}	U^{33}	U^{12}	U^{13}	U^{23}
B1	0.025 (3)	0.031 (3)	0.028 (3)	0.005 (2)	0.006 (2)	-0.001 (2)
C1	0.038 (3)	0.022 (2)	0.020 (2)	0.0017 (18)	-0.0020 (18)	0.0005 (17)
C2	0.036 (3)	0.033 (2)	0.028 (2)	0.0073 (19)	-0.0084 (19)	0.0016 (19)
C3	0.021 (2)	0.026 (2)	0.037 (3)	0.0045 (17)	0.0015 (18)	0.0056 (19)
C4	0.033 (2)	0.026 (2)	0.024 (2)	-0.0042 (18)	0.0010 (18)	0.0035 (18)
C5	0.034 (3)	0.036 (3)	0.032 (2)	-0.0045 (19)	0.0105 (19)	0.009 (2)
C6	0.025 (2)	0.047 (3)	0.023 (2)	-0.0026 (19)	0.0100 (17)	0.0058 (19)
C7	0.028 (2)	0.025 (2)	0.024 (2)	-0.0035 (18)	0.0006 (17)	0.0000 (17)
C8	0.037 (3)	0.029 (2)	0.033 (3)	-0.0087 (19)	-0.0037 (19)	-0.004 (2)
C9	0.048 (3)	0.022 (2)	0.029 (2)	0.0042 (19)	0.004 (2)	-0.0041 (19)
C10	0.017 (2)	0.023 (2)	0.018 (2)	-0.0013 (16)	0.0032 (15)	0.0009 (16)
C11	0.020 (2)	0.025 (2)	0.028 (2)	0.0013 (16)	0.0048 (17)	0.0001 (18)
C12	0.035 (3)	0.019 (2)	0.036 (2)	-0.0033 (18)	0.0028 (19)	-0.0010 (18)
C13	0.027 (2)	0.030 (2)	0.027 (2)	-0.0109 (18)	0.0061 (17)	-0.0006 (18)
C14	0.021 (2)	0.032 (2)	0.032 (2)	0.0007 (17)	0.0050 (17)	-0.0034 (19)
C15	0.027 (2)	0.020 (2)	0.025 (2)	-0.0008 (17)	0.0032 (17)	-0.0006 (17)
C16	0.024 (2)	0.024 (2)	0.018 (2)	0.0070 (16)	0.0031 (16)	-0.0004 (16)
C17	0.025 (2)	0.031 (2)	0.026 (2)	0.0006 (17)	0.0016 (17)	0.0015 (19)
C18	0.029 (2)	0.046 (3)	0.028 (2)	0.004 (2)	-0.0019 (18)	-0.005 (2)
C19	0.035 (3)	0.043 (3)	0.020 (2)	0.015 (2)	0.0022 (19)	0.004 (2)
C20	0.041 (3)	0.032 (2)	0.024 (2)	0.007 (2)	0.0088 (19)	0.005 (2)
C21	0.031 (2)	0.028 (2)	0.022 (2)	-0.0023 (17)	0.0015 (17)	0.0000 (17)
C22	0.020 (2)	0.0188 (19)	0.023 (2)	-0.0035 (16)	0.0047 (16)	-0.0039 (17)
C23	0.027 (2)	0.035 (2)	0.029 (2)	0.0039 (19)	0.0049 (18)	0.0014 (19)
C24	0.035 (3)	0.046 (3)	0.035 (3)	0.009 (2)	0.016 (2)	0.001 (2)
C25	0.021 (2)	0.040 (3)	0.042 (3)	0.0084 (18)	0.004 (2)	-0.006 (2)
C26	0.027 (2)	0.026 (2)	0.030 (2)	0.0022 (17)	-0.0036 (18)	-0.0047 (18)
C27	0.023 (2)	0.024 (2)	0.023 (2)	-0.0002 (16)	0.0053 (16)	-0.0001 (17)
C28	0.021 (2)	0.023 (2)	0.020 (2)	-0.0022 (16)	0.0010 (15)	-0.0024 (17)
C29	0.019 (2)	0.023 (2)	0.034 (2)	-0.0022 (16)	0.0034 (18)	-0.0101 (18)
C30	0.025 (2)	0.050 (3)	0.028 (2)	-0.003 (2)	0.0047 (18)	-0.013 (2)
C31	0.030 (3)	0.068 (3)	0.041 (3)	-0.004 (2)	0.015 (2)	-0.025 (3)

C32	0.022 (3)	0.053 (3)	0.062 (3)	0.004 (2)	0.002 (2)	-0.023 (3)
C33	0.031 (3)	0.043 (3)	0.059 (3)	0.004 (2)	-0.007 (2)	-0.004 (2)
C34	0.025 (3)	0.032 (2)	0.043 (3)	0.0008 (18)	0.0023 (19)	0.005 (2)
N1	0.0232 (18)	0.0173 (16)	0.0244 (18)	-0.0015 (13)	0.0042 (14)	0.0000 (14)
N2	0.0193 (18)	0.0260 (17)	0.0255 (18)	0.0024 (14)	0.0047 (14)	0.0024 (15)
N3	0.0207 (18)	0.0228 (17)	0.0195 (17)	-0.0011 (13)	0.0020 (13)	0.0022 (14)
N4	0.0198 (17)	0.0311 (19)	0.0218 (17)	0.0011 (14)	0.0043 (13)	0.0026 (15)
N5	0.0238 (19)	0.0212 (17)	0.0221 (17)	0.0006 (14)	0.0032 (14)	-0.0026 (14)
N6	0.0270 (19)	0.0232 (18)	0.0257 (19)	0.0004 (14)	0.0067 (14)	-0.0014 (14)
P1	0.0186 (5)	0.0182 (5)	0.0203 (5)	-0.0001 (4)	0.0030 (4)	0.0008 (4)
Ru1	0.01702 (17)	0.01847 (15)	0.01893 (17)	-0.00125 (13)	0.00328 (11)	0.00074 (14)
S1	0.0213 (5)	0.0233 (5)	0.0239 (5)	-0.0014 (4)	0.0042 (4)	0.0021 (4)
S2	0.0216 (6)	0.0252 (5)	0.0225 (5)	-0.0010 (4)	0.0016 (4)	0.0013 (4)

Geometric parameters (Å, °)

B1—N2	1.533 (5)	C18—C19	1.361 (5)
B1—N4	1.535 (5)	C18—H18	0.9500
B1—N6	1.538 (5)	C19—C20	1.378 (5)
B1—H1'	1.0000	C19—H19	0.9500
C1—N1	1.335 (4)	C20—C21	1.391 (5)
C1—C2	1.376 (5)	C20—H20	0.9500
C1—H1	0.9500	C21—H21	0.9500
C2—C3	1.363 (5)	C22—C23	1.386 (5)
C2—H2	0.9500	C22—C27	1.398 (5)
C3—N2	1.342 (4)	C22—P1	1.829 (4)
C3—H3	0.9500	C23—C24	1.386 (5)
C4—N3	1.325 (4)	C23—H23	0.9500
C4—C5	1.382 (5)	C24—C25	1.375 (5)
C4—H4	0.9500	C24—H24	0.9500
C5—C6	1.368 (5)	C25—C26	1.375 (5)
C5—H5	0.9500	C25—H25	0.9500
C6—N4	1.347 (4)	C26—C27	1.378 (5)
C6—H6	0.9500	C26—H26	0.9500
C7—N5	1.334 (4)	C27—H27	0.9500
C7—C8	1.386 (5)	C28—C29	1.461 (5)
C7—H7	0.9500	C28—S2	1.684 (4)
C8—C9	1.355 (5)	C28—S1	1.694 (4)
C8—H8	0.9500	C29—C30	1.392 (5)
C9—N6	1.347 (5)	C29—C34	1.397 (5)
C9—H9	0.9500	C30—C31	1.388 (5)
C10—C11	1.384 (5)	C30—H30	0.9500
C10—C15	1.389 (5)	C31—C32	1.376 (6)
C10—P1	1.834 (3)	C31—H31	0.9500
C11—C12	1.383 (5)	C32—C33	1.364 (6)
C11—H11	0.9500	C32—H32	0.9500
C12—C13	1.376 (5)	C33—C34	1.384 (5)
C12—H12	0.9500	C33—H33	0.9500

C13—C14	1.379 (5)	C34—H34	0.9500
C13—H13	0.9500	N1—N2	1.372 (4)
C14—C15	1.386 (5)	N1—Ru1	2.141 (3)
C14—H14	0.9500	N3—N4	1.366 (4)
C15—H15	0.9500	N3—Ru1	2.098 (3)
C16—C21	1.387 (5)	N5—N6	1.357 (4)
C16—C17	1.389 (5)	N5—Ru1	2.134 (3)
C16—P1	1.844 (4)	P1—Ru1	2.3100 (10)
C17—C18	1.384 (5)	Ru1—S2	2.3213 (10)
C17—H17	0.9500	Ru1—S1	2.3962 (10)
N2—B1—N4	108.0 (3)	C25—C24—C23	120.6 (4)
N2—B1—N6	107.8 (3)	C25—C24—H24	119.7
N4—B1—N6	108.2 (3)	C23—C24—H24	119.7
N2—B1—H1'	110.9	C26—C25—C24	119.5 (4)
N4—B1—H1'	110.9	C26—C25—H25	120.3
N6—B1—H1'	110.9	C24—C25—H25	120.3
N1—C1—C2	111.5 (3)	C25—C26—C27	120.4 (4)
N1—C1—H1	124.2	C25—C26—H26	119.8
C2—C1—H1	124.2	C27—C26—H26	119.8
C3—C2—C1	104.9 (3)	C26—C27—C22	120.8 (3)
C3—C2—H2	127.5	C26—C27—H27	119.6
C1—C2—H2	127.5	C22—C27—H27	119.6
N2—C3—C2	108.7 (3)	C29—C28—S2	124.0 (3)
N2—C3—H3	125.7	C29—C28—S1	126.4 (3)
C2—C3—H3	125.7	S2—C28—S1	109.6 (2)
N3—C4—C5	110.9 (4)	C30—C29—C34	118.5 (4)
N3—C4—H4	124.5	C30—C29—C28	121.0 (4)
C5—C4—H4	124.5	C34—C29—C28	120.6 (3)
C6—C5—C4	105.1 (3)	C31—C30—C29	120.3 (4)
C6—C5—H5	127.5	C31—C30—H30	119.8
C4—C5—H5	127.5	C29—C30—H30	119.8
N4—C6—C5	108.4 (4)	C32—C31—C30	120.1 (4)
N4—C6—H6	125.8	C32—C31—H31	119.9
C5—C6—H6	125.8	C30—C31—H31	119.9
N5—C7—C8	110.2 (3)	C33—C32—C31	120.3 (4)
N5—C7—H7	124.9	C33—C32—H32	119.9
C8—C7—H7	124.9	C31—C32—H32	119.9
C9—C8—C7	105.3 (3)	C32—C33—C34	120.4 (4)
C9—C8—H8	127.4	C32—C33—H33	119.8
C7—C8—H8	127.4	C34—C33—H33	119.8
N6—C9—C8	108.9 (4)	C33—C34—C29	120.4 (4)
N6—C9—H9	125.6	C33—C34—H34	119.8
C8—C9—H9	125.6	C29—C34—H34	119.8
C11—C10—C15	119.1 (3)	C1—N1—N2	105.1 (3)
C11—C10—P1	123.9 (3)	C1—N1—Ru1	137.7 (2)
C15—C10—P1	116.7 (3)	N2—N1—Ru1	117.1 (2)
C12—C11—C10	120.5 (3)	C3—N2—N1	109.7 (3)

C12—C11—H11	119.7	C3—N2—B1	128.4 (3)
C10—C11—H11	119.7	N1—N2—B1	121.8 (3)
C13—C12—C11	120.2 (4)	C4—N3—N4	106.3 (3)
C13—C12—H12	119.9	C4—N3—Ru1	134.1 (3)
C11—C12—H12	119.9	N4—N3—Ru1	119.3 (2)
C12—C13—C14	119.7 (3)	C6—N4—N3	109.3 (3)
C12—C13—H13	120.1	C6—N4—B1	130.3 (3)
C14—C13—H13	120.1	N3—N4—B1	120.4 (3)
C13—C14—C15	120.4 (4)	C7—N5—N6	106.5 (3)
C13—C14—H14	119.8	C7—N5—Ru1	133.3 (3)
C15—C14—H14	119.8	N6—N5—Ru1	120.1 (2)
C14—C15—C10	120.0 (3)	C9—N6—N5	109.2 (3)
C14—C15—H15	120.0	C9—N6—B1	131.7 (3)
C10—C15—H15	120.0	N5—N6—B1	119.1 (3)
C21—C16—C17	117.3 (3)	C22—P1—C10	104.42 (16)
C21—C16—P1	120.7 (3)	C22—P1—C16	102.15 (16)
C17—C16—P1	122.0 (3)	C10—P1—C16	99.36 (16)
C18—C17—C16	121.4 (4)	C22—P1—Ru1	114.77 (11)
C18—C17—H17	119.3	C10—P1—Ru1	116.19 (12)
C16—C17—H17	119.3	C16—P1—Ru1	117.70 (12)
C19—C18—C17	120.0 (4)	N3—Ru1—N5	85.49 (11)
C19—C18—H18	120.0	N3—Ru1—N1	85.83 (11)
C17—C18—H18	120.0	N5—Ru1—N1	84.61 (11)
C18—C19—C20	120.7 (4)	N3—Ru1—P1	96.44 (8)
C18—C19—H19	119.7	N5—Ru1—P1	177.43 (8)
C20—C19—H19	119.7	N1—Ru1—P1	93.82 (8)
C19—C20—C21	118.9 (4)	N3—Ru1—S2	95.15 (8)
C19—C20—H20	120.5	N5—Ru1—S2	88.31 (8)
C21—C20—H20	120.5	N1—Ru1—S2	172.76 (8)
C16—C21—C20	121.7 (4)	P1—Ru1—S2	93.20 (3)
C16—C21—H21	119.2	N3—Ru1—S1	166.07 (8)
C20—C21—H21	119.2	N5—Ru1—S1	89.71 (8)
C23—C22—C27	118.1 (3)	N1—Ru1—S1	106.76 (8)
C23—C22—P1	122.6 (3)	P1—Ru1—S1	88.79 (3)
C27—C22—P1	118.8 (3)	S2—Ru1—S1	71.61 (3)
C22—C23—C24	120.5 (4)	C28—S1—Ru1	88.02 (12)
C22—C23—H23	119.7	C28—S2—Ru1	90.79 (12)
C24—C23—H23	119.7		
