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Bis(2,2′-bipyridine-κ^2N,N')dichloridoplatinum(IV) dichloride monohydrate

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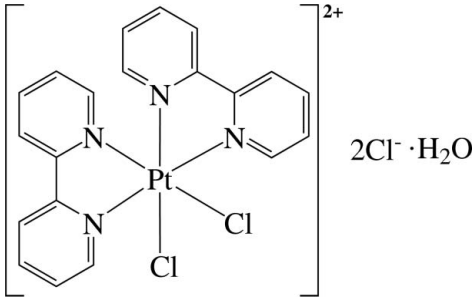
Key indicators: single-crystal X-ray study; T = 293 K; mean σ(C–C) = 0.005 Å; R factor = 0.016; wR factor = 0.038; data-to-parameter ratio = 16.5.

In the title complex, $[PtCl_2(C_{10}H_8N_2)_2]Cl_2\cdot H_2O$, the Pt^{4+} ion is six-coordinated in a distorted octahedral environment by four N atoms from the two 2,2′-bipyridine ligands and two Cl atoms. As a result of the different *trans* influences of the N and Cl atoms, the Pt—N bonds *trans* to the Cl atom are slightly longer than those *trans* to the N atom. The compound displays intermolecular hydrogen bonding between the water molecule and the Cl anions. There are intermolecular π–π interactions between adjacent pyridine rings, with a centroid–centroid distance of 3.962 Å.

Related literature

For related literature, see: Hambley (1986); Hojjat Kashani *et al.* (2008).



Experimental

Crystal data

$[PtCl_2(C_{10}H_8N_2)_2]Cl_2\cdot H_2O$
M_r = 667.27
Orthorhombic, $P2_12_12_1$
a = 11.1345 (12) Å
b = 11.5867 (12) Å
c = 17.0873 (19) Å
V = 2204.5 (4) Å^3
Z = 4
Mo $K\alpha$ radiation
μ = 6.87 mm^{-1}
T = 293 (2) K
0.35 × 0.20 × 0.15 mm

Data collection

Bruker SMART 1000 CCD diffractometer
Absorption correction: multi-scan (*SADABS*; Bruker, 2000)
T_{min} = 0.251, T_{max} = 0.357
12649 measured reflections
4462 independent reflections
4284 reflections with $I > 2\sigma(I)$
R_{int} = 0.017

Refinement

$R[F^2 > 2\sigma(F^2)]$ = 0.016
$wR(F^2)$ = 0.038
S = 0.84
4462 reflections
271 parameters
H-atom parameters constrained
$\Delta\rho_{max}$ = 0.95 e Å^{-3}
$\Delta\rho_{min}$ = −0.53 e Å^{-3}
Absolute structure: Flack (1983), 1901 Friedel pairs
Flack parameter: −0.006 (4)

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

D—H···A	D—H	H···A	D···A	D—H···A
O1—H1A···Cl3^{i}	1.033	2.21	3.150 (3)	149.79 (16)
O1—H1B···Cl4^{ii}	0.924	2.31	3.139 (3)	149.3 (2)

Symmetry codes: (i) $x-\frac{1}{2}, -y+\frac{1}{2}, -z+1$; (ii) $-x, y-\frac{1}{2}, -z+\frac{1}{2}$.

Data collection: *SMART* (Bruker, 2000); cell refinement: *SAINT* (Bruker, 2000); data reduction: *SAINT*; program(s) used to solve structure: *SHELXS97* (Sheldrick, 2008); program(s) used to refine structure: *SHELXL97* (Sheldrick, 2008); molecular graphics: *ORTEP-3* (Farrugia, 1997); software used to prepare material for publication: *SHELXL97*.

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Supplementary data and figures for this paper are available from the IUCr electronic archives (Reference: BT2846).

References

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supplementary materials

Acta Cryst. (2009). E**65**, m180 [**doi:10.1107/S1600536809000725**]

Bis(2,2'-bipyridine-κ^2N,N')dichloridoplatinum(IV) dichloride monohydrate

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Comment

In the title complex, $[PtCl_2(C_{10}H_8N_2)_2]Cl_2.H_2O$, the central Pt^{4+} ion is six-coordinated in a distorted octahedral environment by four N atoms from the two 2,2'-bipyridine ligands and two Cl atoms (Fig. 1). The main contributions to the distortion are the tight N—Pt—N chelate angles (80.33 (10)° and 80.30 (10)°), which result in non-linear *trans* axes (<Cl1—Pt1—N1 = 176.73 (7)°, <Cl2—Pt1—N4 = 176.91 (7)° and <N2—Pt1—N3 = 176.52 (10)°).

Because of the different *trans* influences of the N and Cl atoms, the Pt—N bonds *trans* to the Cl atom (lengths: 2.040 (2) and 2.037 (3) Å) are slightly longer than those *trans* to the N atom (lengths: 2.029 (2) and 2.028 (2) Å).

The compound displays intermolecular hydrogen bonding between the solvent H_2O molecule and the Cl anions (Table 1). There is also an intermolecular π-π interaction between the pyridine ring containing N1 and the one containg N3 at 1/2+x,1/2-y,-z, with a centroid-centroid distance of 3.962 Å and with a dihedral angle between the ring planes of 20.3°.

Experimental

To a solution of K_2PtCl_6 (0.3068 g, 0.631 mmol) in H_2O (20 ml) was added 2,2'-bipyridine (0.0971 g, 0.622 mmol) in MeOH (10 ml), and stirred for 2 h under heating. The formed precipitate was separated by filtration and washed with water and MeOH and dried under vacuum, to give a yellow powder (0.1185 g). Crystals suitable for X-ray analysis were obtained by slow evaporation from a CH_2Cl_2 solution.

Refinement

H atoms were positioned geometrically and allowed to ride on their respective parent atoms [C—H = 0.93 Å and U_{iso}(H) = 1.2U_{eq}(C)]. The H atoms of the solvent H_2O molecule were located from Fourier difference maps, but not refined.

Figures

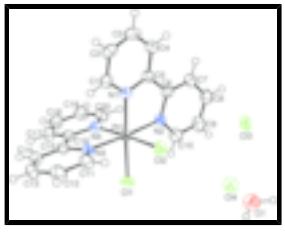


Fig. 1. The structure of the title compound, with displacement ellipsoids drawn at the 50% probability level for non-H atoms.

Bis(2,2'-bipyridine-κ^2*N*,*N*')dichloridoplatinum(IV) dichloride monohydrate

Crystal data

$[PtCl_2(C_{10}H_8N_2)_2]Cl_2·H_2O$	F_{000} = 1280
M_r = 667.27	D_x = 2.011 Mg m^{-3}
Orthorhombic, $P2_12_12_1$	Mo $K\alpha$ radiation λ = 0.71073 Å
Hall symbol: P 2ac 2ab	Cell parameters from 958 reflections
a = 11.1345 (12) Å	θ = 2.4–26.4º
b = 11.5867 (12) Å	μ = 6.87 mm^{-1}
c = 17.0873 (19) Å	T = 293 (2) K
V = 2204.5 (4) Å^3	Stick, colorless
Z = 4	0.35 × 0.20 × 0.15 mm

Data collection

Bruker SMART 1000 CCD diffractometer	4462 independent reflections
Radiation source: fine-focus sealed tube	4284 reflections with $I > 2\sigma(I)$
Monochromator: graphite	R_{int} = 0.017
T = 293(2) K	θ_{max} = 26.4º
φ and ω scans	θ_{min} = 2.2º
Absorption correction: Multi-scan (SADABS; Bruker, 2000)	h = −12→13
T_{min} = 0.251, T_{max} = 0.357	k = −14→14
12649 measured reflections	l = −21→16

Refinement

Refinement on F^2	Hydrogen site location: inferred from neighbouring sites
Least-squares matrix: full	H-atom parameters constrained
$R[F^2 > 2\sigma(F^2)]$ = 0.016	$w = 1/[\sigma^2(F_o^2)]$ where $P = (F_o^2 + 2F_c^2)/3$
$wR(F^2)$ = 0.038	$(\Delta/\sigma)_{max}$ = 0.003
S = 0.84	$\Delta\rho_{max}$ = 0.95 e Å^{-3}
4462 reflections	$\Delta\rho_{min}$ = −0.53 e Å^{-3}
271 parameters	Extinction correction: none
Primary atom site location: structure-invariant direct methods	Absolute structure: Flack (1983), 1901 Friedel pairs
Secondary atom site location: difference Fourier map	Flack parameter: −0.006 (4)

Special details

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 (Å^2)

	x	*y*	*z*	U_{iso}*/U_{eq}
Pt1	0.520023 (9)	0.241999 (9)	0.124231 (6)	0.02622 (4)
Cl1	0.35373 (7)	0.12668 (7)	0.10471 (5)	0.0410 (2)
Cl2	0.40172 (7)	0.40216 (7)	0.14570 (5)	0.0394 (2)
Cl3	0.36627 (9)	0.07710 (8)	0.54638 (6)	0.0517 (2)
Cl4	0.15636 (8)	0.59942 (8)	0.20502 (5)	0.0421 (2)
N1	0.6674 (2)	0.3407 (2)	0.14792 (15)	0.0272 (6)
N2	0.5239 (2)	0.2274 (2)	0.24255 (14)	0.0290 (5)
N3	0.5272 (2)	0.2554 (2)	0.00599 (14)	0.0288 (5)
N4	0.6210 (2)	0.1002 (2)	0.09926 (15)	0.0309 (6)
C1	0.7301 (3)	0.3999 (3)	0.0943 (2)	0.0340 (7)
H1	0.7110	0.3925	0.0416	0.041*
C2	0.8225 (3)	0.4716 (3)	0.1167 (2)	0.0389 (8)
H2	0.8644	0.5140	0.0794	0.047*
C3	0.8524 (3)	0.4801 (3)	0.1945 (2)	0.0413 (8)
H3	0.9159	0.5270	0.2099	0.050*
C4	0.7876 (3)	0.4185 (3)	0.2500 (2)	0.0364 (8)
H4	0.8078	0.4227	0.3027	0.044*
C5	0.6933 (3)	0.3516 (3)	0.22561 (18)	0.0284 (7)
C6	0.6124 (3)	0.2891 (3)	0.27847 (18)	0.0282 (7)
C7	0.6188 (3)	0.2940 (3)	0.35849 (19)	0.0391 (8)
H7	0.6792	0.3364	0.3827	0.047*
C8	0.5347 (3)	0.2355 (3)	0.4032 (2)	0.0415 (8)
H8	0.5378	0.2380	0.4575	0.050*
C9	0.4458 (3)	0.1728 (3)	0.3651 (2)	0.0452 (9)
H9	0.3886	0.1327	0.3940	0.054*
C10	0.4424 (3)	0.1702 (3)	0.2852 (2)	0.0399 (8)
H10	0.3827	0.1281	0.2601	0.048*
C11	0.4779 (3)	0.3413 (3)	−0.0360 (2)	0.0386 (8)
H11	0.4407	0.4024	−0.0103	0.046*
C12	0.4818 (3)	0.3402 (3)	−0.1165 (2)	0.0453 (9)
H12	0.4465	0.3996	−0.1450	0.054*
C13	0.5380 (3)	0.2508 (3)	−0.15455 (19)	0.0413 (8)
H13	0.5409	0.2489	−0.2089	0.050*

C14	0.5907 (3)	0.1629 (3)	−0.11082 (19)	0.0371 (8)
H14	0.6302	0.1024	−0.1356	0.045*
C15	0.5836 (3)	0.1666 (3)	−0.03080 (19)	0.0294 (7)
C16	0.6334 (3)	0.0775 (3)	0.02149 (18)	0.0299 (7)
C17	0.6884 (3)	−0.0219 (3)	−0.0035 (2)	0.0404 (8)
H17	0.6942	−0.0385	−0.0566	0.048*
C18	0.7351 (3)	−0.0971 (3)	0.0517 (2)	0.0444 (9)
H18	0.7734	−0.1644	0.0357	0.053*
C19	0.7249 (3)	−0.0725 (3)	0.1296 (2)	0.0435 (8)
H19	0.7567	−0.1223	0.1669	0.052*
C20	0.6665 (3)	0.0280 (3)	0.1524 (2)	0.0366 (8)
H20	0.6589	0.0451	0.2054	0.044*
O1	0.0454 (3)	0.2475 (2)	0.37482 (18)	0.0807 (11)
H1A	0.0141	0.3103	0.4125	0.080*
H1B	−0.0004	0.1831	0.3636	0.080*

Atomic displacement parameters (Å^2)

	U^{11}	U^{22}	U^{33}	U^{12}	U^{13}	U^{23}
Pt1	0.02736 (6)	0.02822 (6)	0.02310 (6)	−0.00141 (5)	−0.00187 (4)	0.00024 (5)
Cl1	0.0367 (4)	0.0447 (5)	0.0416 (5)	−0.0117 (4)	−0.0046 (3)	−0.0014 (4)
Cl2	0.0400 (4)	0.0375 (4)	0.0408 (5)	0.0082 (4)	−0.0011 (3)	−0.0036 (3)
Cl3	0.0656 (6)	0.0536 (6)	0.0360 (5)	−0.0087 (5)	0.0145 (4)	−0.0081 (4)
Cl4	0.0430 (5)	0.0454 (5)	0.0379 (5)	0.0001 (4)	−0.0017 (4)	−0.0040 (4)
N1	0.0243 (13)	0.0257 (13)	0.0314 (16)	−0.0014 (10)	−0.0031 (11)	0.0011 (10)
N2	0.0319 (13)	0.0301 (13)	0.0251 (13)	−0.0025 (14)	0.0013 (10)	0.0010 (10)
N3	0.0294 (13)	0.0332 (13)	0.0238 (12)	0.0022 (16)	−0.0056 (9)	0.0013 (10)
N4	0.0294 (14)	0.0309 (14)	0.0325 (15)	−0.0022 (11)	−0.0021 (11)	−0.0006 (11)
C1	0.0361 (18)	0.0362 (18)	0.0297 (18)	0.0021 (15)	0.0030 (14)	0.0024 (14)
C2	0.0323 (17)	0.0387 (18)	0.046 (2)	−0.0046 (13)	0.0074 (17)	0.0082 (17)
C3	0.0320 (18)	0.0377 (19)	0.054 (2)	−0.0076 (15)	−0.0029 (16)	−0.0009 (16)
C4	0.0338 (18)	0.042 (2)	0.033 (2)	−0.0010 (15)	−0.0042 (14)	−0.0032 (15)
C5	0.0288 (16)	0.0294 (16)	0.0271 (17)	0.0042 (13)	−0.0025 (13)	0.0009 (13)
C6	0.0277 (16)	0.0301 (15)	0.0269 (16)	0.0028 (12)	−0.0013 (12)	−0.0002 (12)
C7	0.0398 (19)	0.0493 (19)	0.0281 (19)	−0.0004 (15)	−0.0020 (14)	−0.0022 (15)
C8	0.050 (2)	0.050 (2)	0.0240 (16)	0.003 (2)	0.0016 (13)	0.0056 (14)
C9	0.058 (2)	0.0409 (19)	0.037 (2)	−0.0088 (16)	0.0134 (18)	0.0074 (16)
C10	0.047 (2)	0.0370 (18)	0.035 (2)	−0.0099 (16)	0.0033 (16)	0.0032 (14)
C11	0.044 (2)	0.0380 (17)	0.0336 (19)	0.0054 (17)	−0.0062 (16)	0.0047 (14)
C12	0.051 (2)	0.0494 (19)	0.035 (2)	0.0019 (17)	−0.0079 (18)	0.0104 (16)
C13	0.0481 (19)	0.052 (2)	0.0233 (15)	0.000 (3)	−0.0028 (13)	0.0036 (15)
C14	0.0406 (19)	0.0441 (18)	0.0267 (19)	−0.0010 (15)	0.0029 (15)	−0.0038 (14)
C15	0.0268 (16)	0.0338 (17)	0.0277 (17)	−0.0020 (13)	−0.0035 (13)	0.0011 (13)
C16	0.0288 (17)	0.0327 (17)	0.0283 (18)	−0.0023 (13)	−0.0017 (13)	−0.0008 (13)
C17	0.045 (2)	0.039 (2)	0.037 (2)	0.0053 (17)	0.0047 (15)	−0.0058 (15)
C18	0.044 (2)	0.038 (2)	0.051 (3)	0.0120 (17)	0.0043 (17)	−0.0011 (17)
C19	0.0421 (19)	0.0395 (18)	0.049 (2)	0.0079 (15)	−0.0068 (18)	0.0024 (19)
C20	0.0423 (19)	0.0386 (18)	0.0287 (18)	−0.0012 (15)	−0.0087 (15)	0.0029 (14)

O1	0.0528 (17)	0.084 (2)	0.106 (3)	−0.0042 (15)	−0.0045 (15)	−0.043 (2)

Geometric parameters (Å, °)

Pt1—N3	2.028 (2)	C7—H7	0.9300
Pt1—N2	2.029 (2)	C8—C9	1.389 (5)
Pt1—N4	2.037 (3)	C8—H8	0.9300
Pt1—N1	2.040 (2)	C9—C10	1.366 (5)
Pt1—Cl2	2.3051 (8)	C9—H9	0.9300
Pt1—Cl1	2.3076 (8)	C10—H10	0.9300
N1—C1	1.341 (4)	C11—C12	1.377 (5)
N1—C5	1.364 (4)	C11—H11	0.9300
N2—C10	1.339 (4)	C12—C13	1.375 (5)
N2—C6	1.364 (4)	C12—H12	0.9300
N3—C11	1.344 (4)	C13—C14	1.393 (5)
N3—C15	1.360 (4)	C13—H13	0.9300
N4—C20	1.335 (4)	C14—C15	1.370 (4)
N4—C16	1.361 (4)	C14—H14	0.9300
C1—C2	1.377 (4)	C15—C16	1.474 (4)
C1—H1	0.9300	C16—C17	1.373 (4)
C2—C3	1.373 (5)	C17—C18	1.385 (5)
C2—H2	0.9300	C17—H17	0.9300
C3—C4	1.389 (5)	C18—C19	1.367 (5)
C3—H3	0.9300	C18—H18	0.9300
C4—C5	1.370 (4)	C19—C20	1.389 (4)
C4—H4	0.9300	C19—H19	0.9300
C5—C6	1.467 (4)	C20—H20	0.9300
C6—C7	1.370 (4)	O1—H1A	1.033
C7—C8	1.385 (5)	O1—H1B	0.924
N3—Pt1—N2	176.52 (10)	C7—C6—C5	124.2 (3)
N3—Pt1—N4	80.30 (10)	C6—C7—C8	119.6 (3)
N2—Pt1—N4	97.47 (10)	C6—C7—H7	120.2
N3—Pt1—N1	97.07 (10)	C8—C7—H7	120.2
N2—Pt1—N1	80.33 (10)	C7—C8—C9	118.7 (3)
N4—Pt1—N1	92.84 (9)	C7—C8—H8	120.7
N3—Pt1—Cl2	96.85 (7)	C9—C8—H8	120.7
N2—Pt1—Cl2	85.44 (7)	C10—C9—C8	119.9 (3)
N4—Pt1—Cl2	176.91 (7)	C10—C9—H9	120.1
N1—Pt1—Cl2	88.68 (7)	C8—C9—H9	120.1
N3—Pt1—Cl1	86.10 (7)	N2—C10—C9	121.0 (3)
N2—Pt1—Cl1	96.47 (7)	N2—C10—H10	119.5
N4—Pt1—Cl1	86.88 (7)	C9—C10—H10	119.5
N1—Pt1—Cl1	176.73 (7)	N3—C11—C12	120.9 (3)
Cl2—Pt1—Cl1	91.76 (3)	N3—C11—H11	119.5
C1—N1—C5	120.5 (3)	C12—C11—H11	119.5
C1—N1—Pt1	124.8 (2)	C13—C12—C11	119.6 (3)
C5—N1—Pt1	114.53 (19)	C13—C12—H12	120.2
C10—N2—C6	120.3 (3)	C11—C12—H12	120.2
C10—N2—Pt1	124.7 (2)	C12—C13—C14	119.3 (3)

C6—N2—Pt1	114.83 (19)	C12—C13—H13	120.3
C11—N3—C15	120.2 (3)	C14—C13—H13	120.3
C11—N3—Pt1	124.9 (2)	C15—C14—C13	119.2 (3)
C15—N3—Pt1	114.88 (19)	C15—C14—H14	120.4
C20—N4—C16	120.3 (3)	C13—C14—H14	120.4
C20—N4—Pt1	124.9 (2)	N3—C15—C14	120.8 (3)
C16—N4—Pt1	114.6 (2)	N3—C15—C16	115.1 (3)
N1—C1—C2	120.6 (3)	C14—C15—C16	124.1 (3)
N1—C1—H1	119.7	N4—C16—C17	120.7 (3)
C2—C1—H1	119.7	N4—C16—C15	114.8 (3)
C3—C2—C1	119.5 (3)	C17—C16—C15	124.5 (3)
C3—C2—H2	120.2	C16—C17—C18	118.9 (3)
C1—C2—H2	120.2	C16—C17—H17	120.6
C2—C3—C4	119.9 (3)	C18—C17—H17	120.6
C2—C3—H3	120.0	C19—C18—C17	120.1 (3)
C4—C3—H3	120.0	C19—C18—H18	119.9
C5—C4—C3	118.7 (3)	C17—C18—H18	119.9
C5—C4—H4	120.6	C18—C19—C20	119.1 (3)
C3—C4—H4	120.6	C18—C19—H19	120.4
N1—C5—C4	120.7 (3)	C20—C19—H19	120.4
N1—C5—C6	115.0 (3)	N4—C20—C19	120.8 (3)
C4—C5—C6	124.2 (3)	N4—C20—H20	119.6
N2—C6—C7	120.6 (3)	C19—C20—H20	119.6
N2—C6—C5	115.2 (3)	H1A—O1—H1B	120.8
N3—Pt1—N1—C1	6.2 (3)	C3—C4—C5—C6	−175.4 (3)
N2—Pt1—N1—C1	−176.1 (3)	C10—N2—C6—C7	−0.4 (4)
N4—Pt1—N1—C1	86.8 (3)	Pt1—N2—C6—C7	174.8 (2)
Cl2—Pt1—N1—C1	−90.5 (2)	C10—N2—C6—C5	−177.7 (3)
N3—Pt1—N1—C5	−179.5 (2)	Pt1—N2—C6—C5	−2.6 (3)
N2—Pt1—N1—C5	−1.9 (2)	N1—C5—C6—N2	1.0 (4)
N4—Pt1—N1—C5	−99.0 (2)	C4—C5—C6—N2	179.9 (3)
Cl2—Pt1—N1—C5	83.7 (2)	N1—C5—C6—C7	−176.2 (3)
N4—Pt1—N2—C10	−91.0 (3)	C4—C5—C6—C7	2.7 (5)
N1—Pt1—N2—C10	177.4 (3)	N2—C6—C7—C8	0.2 (5)
Cl2—Pt1—N2—C10	87.9 (2)	C5—C6—C7—C8	177.3 (3)
Cl1—Pt1—N2—C10	−3.3 (3)	C6—C7—C8—C9	0.1 (5)
N4—Pt1—N2—C6	94.0 (2)	C7—C8—C9—C10	−0.2 (5)
N1—Pt1—N2—C6	2.4 (2)	C6—N2—C10—C9	0.3 (5)
Cl2—Pt1—N2—C6	−87.0 (2)	Pt1—N2—C10—C9	−174.4 (3)
Cl1—Pt1—N2—C6	−178.28 (19)	C8—C9—C10—N2	0.0 (5)
N4—Pt1—N3—C11	−178.1 (3)	C15—N3—C11—C12	1.1 (5)
N1—Pt1—N3—C11	−86.4 (3)	Pt1—N3—C11—C12	−176.5 (2)
Cl2—Pt1—N3—C11	3.1 (2)	N3—C11—C12—C13	−0.8 (5)
Cl1—Pt1—N3—C11	94.4 (2)	C11—C12—C13—C14	−0.3 (5)
N4—Pt1—N3—C15	4.2 (2)	C12—C13—C14—C15	1.0 (5)
N1—Pt1—N3—C15	95.9 (2)	C11—N3—C15—C14	−0.3 (4)
Cl2—Pt1—N3—C15	−174.61 (19)	Pt1—N3—C15—C14	177.5 (2)
Cl1—Pt1—N3—C15	−83.3 (2)	C11—N3—C15—C16	−179.7 (3)
N3—Pt1—N4—C20	178.5 (3)	Pt1—N3—C15—C16	−1.8 (3)

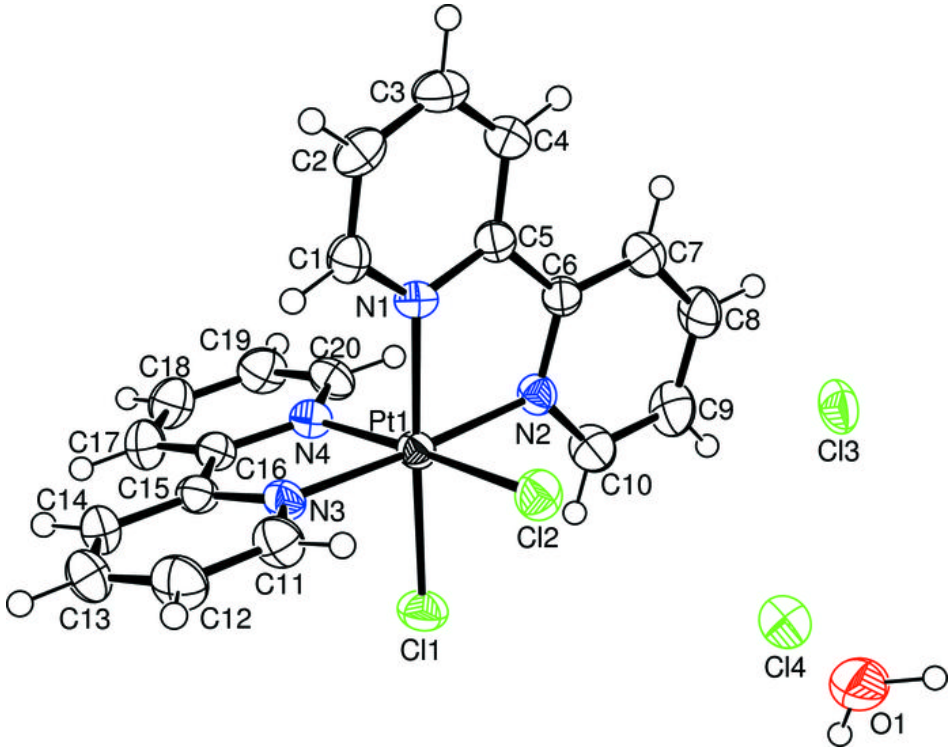
N2—Pt1—N4—C20	1.2 (3)	C13—C14—C15—N3	−0.7 (5)
N1—Pt1—N4—C20	81.8 (3)	C13—C14—C15—C16	178.6 (3)
Cl1—Pt1—N4—C20	−94.9 (3)	C20—N4—C16—C17	2.6 (5)
N3—Pt1—N4—C16	−6.0 (2)	Pt1—N4—C16—C17	−173.2 (2)
N2—Pt1—N4—C16	176.7 (2)	C20—N4—C16—C15	−177.5 (3)
N1—Pt1—N4—C16	−102.7 (2)	Pt1—N4—C16—C15	6.8 (3)
Cl1—Pt1—N4—C16	80.5 (2)	N3—C15—C16—N4	−3.3 (4)
C5—N1—C1—C2	0.8 (5)	C14—C15—C16—N4	177.4 (3)
Pt1—N1—C1—C2	174.7 (2)	N3—C15—C16—C17	176.6 (3)
N1—C1—C2—C3	1.6 (5)	C14—C15—C16—C17	−2.7 (5)
C1—C2—C3—C4	−1.5 (5)	N4—C16—C17—C18	−2.3 (5)
C2—C3—C4—C5	−1.0 (5)	C15—C16—C17—C18	177.8 (3)
C1—N1—C5—C4	−3.4 (4)	C16—C17—C18—C19	0.7 (5)
Pt1—N1—C5—C4	−177.9 (2)	C17—C18—C19—C20	0.6 (5)
C1—N1—C5—C6	175.6 (3)	C16—N4—C20—C19	−1.3 (5)
Pt1—N1—C5—C6	1.1 (3)	Pt1—N4—C20—C19	174.0 (2)
C3—C4—C5—N1	3.4 (5)	C18—C19—C20—N4	−0.3 (5)

Hydrogen-bond geometry (Å, °)

D—H···*A*	*D*—H	H···*A*	*D*···*A*	*D*—H···*A*
O1—H1A···Cl3^{i}	1.033	2.21	3.150 (3)	149.79 (16)
O1—H1B···Cl4^{ii}	0.924	2.31	3.139 (3)	149.3 (2)

Symmetry codes: (i) *x*−1/2, −*y*+1/2, −*z*+1; (ii) −*x*, *y*−1/2, −*z*+1/2.

Fig. 1


C3
C4
C2
C5
C1
C6
C7
N1
C8
C19
C20
C18
Pt1
C9
N2
Cl3
C17
N4
C16
C15
C10
C14
N3
Cl2
C11
C13
C12
Cl1
Cl4
O1