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Tetrakis[1-phenyl-3-(1*H*-1,2,4-triazol-1-yl-κN^4)propan-1-one]bis(thiocyanato-κN)manganese(II)

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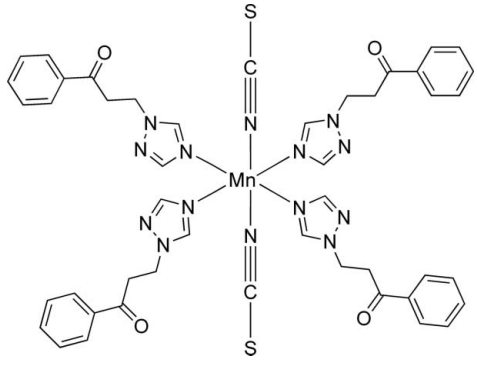
Key indicators: single-crystal X-ray study; T = 293 K; mean σ(C–C) = 0.004 Å; R factor = 0.038; wR factor = 0.089; data-to-parameter ratio = 13.4.

In the mononuclear title complex, $[Mn(NCS)_2(C_{11}H_{11}N_3O)_4]$, the Mn^{II} atom, lying on an inversion center, is coordinated by two monodentate thiocyanate anions and four monodentate 1-phenyl-3-(1*H*-1,2,4-triazol-1-yl)propan-1-one ligands in a distorted octahedral geometry. Each complex molecule is linked to four neighboring ones by weak C—H···N and C—H···S hydrogen bonds, forming a two-dimensional sheet parallel to (001).

Related literature

For general background to self-assembly of supramolecular systems, see: Beatty (2003); Braga *et al.* (2003). For a related structure, see: Guo & Cai (2007).



Experimental

Crystal data

$[Mn(NCS)_2(C_{11}H_{11}N_3O)_4]$ M_r = 976.03
Triclinic, $P\overline{1}$
a = 7.9326 (17) Å
b = 11.845 (3) Å
c = 13.740 (3) Å
α = 69.240 (3)°
β = 75.417 (3)°
γ = 81.686 (3)°
V = 1166.1 (5) Å^3
Z = 1
Mo $K\alpha$ radiation
μ = 0.43 mm^{-1}
T = 293 K
0.20 × 0.18 × 0.14 mm

Data collection

Bruker APEXII CCD diffractometer
Absorption correction: multi-scan (*SADABS*; Sheldrick, 1996)
T_{min} = 0.919, T_{max} = 0.942
6410 measured reflections
4075 independent reflections
2840 reflections with $I > 2\sigma(I)$
R_{int} = 0.021

Refinement

$R[F^2 > 2\sigma(F^2)]$ = 0.038
$wR(F^2)$ = 0.089
S = 1.06
4075 reflections
304 parameters
H-atom parameters constrained
$\Delta\rho_{max}$ = 0.19 e Å^{-3}
$\Delta\rho_{min}$ = −0.23 e Å^{-3}

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

D—H···A	D—H	H···A	D···A	D—H···A
C12—H12···N2^{i}	0.93	2.62	3.436 (3)	146
C18—H18···S1^{ii}	0.93	2.82	3.725 (3)	164

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

Data collection: *APEX2* (Bruker, 2007); cell refinement: *SAINT* (Bruker, 2007); data reduction: *SAINT*; program(s) used to solve structure: *SHELXS97* (Sheldrick, 2008); program(s) used to refine structure: *SHELXL97* (Sheldrick, 2008); molecular graphics: *SHELXTL* (Sheldrick, 2008) and *DIAMOND* (Brandenburg & Berndt, 1999); software used to prepare material for publication: *SHELXTL*.

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Supplementary data and figures for this paper are available from the IUCr electronic archives (Reference: HY2378).

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supplementary materials

Acta Cryst. (2010). E**66**, m1605 [**doi:10.1107/S1600536810047112**]

Tetrakis[1-phenyl-3-(1*H*-1,2,4-triazol-1-yl-κN^4)propan-1-one]bis(thiocyanato-κN)manganese(II)

H. Cai, Y. Guo and J.-G. Li

Comment

Self-assembly processes directed by either hydrogen-bonding interactions or metal coordinations have been extensively utilized in crystal engineering to construct supramolecular systems with novel structures and properties due to their inherent strength and reliability (Braga *et al.*, 2003). Proper selection of metal ions and ligands with suitable functionalized groups is the key issue in designing and self-assembling of coordination supramolecules (Beatty, 2003). Recently, we have initiated a research program of synthesizing supramolecules based on pseudohalide and flexible ligand, which consists of a propanone unit substituted with an imidazole and a phenyl group (Guo & Cai, 2007). To further explore this series, we synthesized the title compound, a new Mn^{II} complex based on the mixed ligands, thiocyanate and 3-(1*H*-1,2,4-triazol-1-yl)-1-phenyl-propan-1-one (*L*).

In the molecular structure (Fig. 1) of the mononuclear title complex, the Mn^{II} atom is six-coordinated by four mono-dentate *L* ligands, forming the equatorial plane and two N atoms from two monodentate NCS^- anions in the axial positions, displaying an MnN_6 octahedral geometry. The triazol and phenyl rings in each of the ligands are not coplanar. The dihedral angels formed by the least-squares planes of the phenyl and triazole rings are 53.8 (2) and 69.6 (2)°. In the crystal, weak intermolecular C—H···N and C—H···S hydrogen bonds (Table 1) connect the complex molecules into a two-dimensional supramolecular sheet parallel to (0 0 1), as shown in Fig. 2.

Experimental

$MnCl_2.4H_2O$ (19.8 mg, 0.1 mmol), 3-(1*H*-1,2,4-triazol-1-yl)-1-phenylpropan-1-one (22.3 mg, 0.1 mmol) and NH_4SCN (7.6 mg, 0.1 mmol) were mixed in a CH_3CN/H_2O (20 ml, v/v 1:1) solution with vigorous stirring for *ca* 30 min. The resulting solution was filtered and left to stand at room temperature. Colorless block crystals of the title compound suitable for X-ray analysis were obtained in a 60% yield by slow evaporation of the solvent over a period of one week. Analysis, calculated for $C_{46}H_{44}MnN_{14}O_4S_2$: C 56.61, H 4.54, N 20.09%; found: C 56.45, H 4.43, N 20.12%.

Refinement

Although all H atoms were visible in difference Fourier maps, they were finally placed in geometrically calculated positions and refined as riding atoms, with C—H = 0.93 (aromatic) and 0.97 (methylene) Å and with $U_{iso}(H) = 1.2U_{eq}(C)$.

Figures

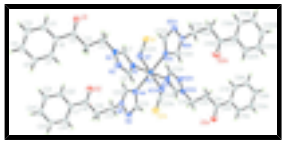


Fig. 1. The molecular structure of the title compound, showing the 30% probability ellipsoids. [Symmetry code: (A) -x, -y, 2-z.]

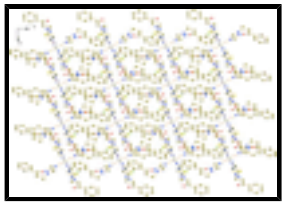


Fig. 2. The two-dimensional sheet structure of the title compound, showing C—H···N and C—H···S hydrogen bongs as red and green dashed lines.

Tetrakis[1-phenyl-3-(1*H*-1,2,4-triazol-1-yl-κ*N*^{4})propan-1- one]bis(thiocyanato-κ*N*)manganese(II)

Crystal data

$[Mn(NCS)_2(C_{11}H_{11}N_3O)_4]$	$Z = 1$
$M_r = 976.03$	$F(000) = 507$
Triclinic, $P\overline{1}$	$D_x = 1.390$ Mg m^{-3}
Hall symbol: -P 1	Mo $K\alpha$ radiation, $\lambda = 0.71073$ Å
$a = 7.9326$ (17) Å	Cell parameters from 1514 reflections
$b = 11.845$ (3) Å	$\theta = 2.8$–$22.4°$
$c = 13.740$ (3) Å	$\mu = 0.43$ mm^{-1}
$\alpha = 69.240$ (3)°	$T = 293$ K
$\beta = 75.417$ (3)°	Block, colorless
$\gamma = 81.686$ (3)°	0.20 × 0.18 × 0.14 mm
$V = 1166.1$ (5) Å^3	

Data collection

Bruker APEXII CCD diffractometer	4075 independent reflections
Radiation source: fine-focus sealed tube	2840 reflections with $I > 2\sigma(I)$
graphite	$R_{int} = 0.021$
φ and ω scans	$\theta_{max} = 25.0°$, $\theta_{min} = 2.0°$
Absorption correction: multi-scan (*SADABS*; Sheldrick, 1996)	$h = -8 \rightarrow 9$
$T_{min} = 0.919$, $T_{max} = 0.942$	$k = -13 \rightarrow 14$
6410 measured reflections	$l = -14 \rightarrow 16$

Refinement

Refinement on F^2	Primary atom site location: structure-invariant direct methods
Least-squares matrix: full	Secondary atom site location: difference Fourier map
$R[F^2 > 2\sigma(F^2)] = 0.038$	Hydrogen site location: inferred from neighbouring sites
$wR(F^2) = 0.089$	H-atom parameters constrained
$S = 1.06$	$w = 1/[\sigma^2(F_o^2) + (0.0394P)^2 + 0.0024P]$ where $P = (F_o^2 + 2F_c^2)/3$
4075 reflections	$(\Delta/\sigma)_{max} < 0.001$
304 parameters	$\Delta\rho_{max} = 0.19$ e Å^{-3}

0 restraints $\Delta\rho_{min} = -0.23$ e Å^{-3}

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

	x	*y*	*z*	U_{iso}*/U_{eq}
Mn1	0.0000	0.0000	1.0000	0.03670 (15)
S1	−0.34152 (10)	−0.18300 (7)	0.86875 (6)	0.0709 (2)
O1	0.7108 (2)	0.12219 (16)	0.41612 (14)	0.0696 (5)
O2	−0.2798 (3)	0.43808 (16)	0.66598 (16)	0.0763 (6)
N1	0.2363 (2)	0.06526 (16)	0.87054 (14)	0.0429 (5)
N2	0.4935 (2)	0.14559 (18)	0.78206 (15)	0.0532 (5)
N3	0.4371 (2)	0.09195 (16)	0.72510 (14)	0.0425 (5)
N4	−0.1390 (2)	0.18782 (15)	0.96560 (14)	0.0413 (5)
N5	−0.2623 (3)	0.36007 (18)	0.99367 (16)	0.0532 (5)
N6	−0.3215 (2)	0.34167 (17)	0.91665 (15)	0.0441 (5)
N7	−0.1117 (3)	−0.03665 (19)	0.88320 (16)	0.0538 (5)
C1	0.2858 (3)	0.0449 (2)	0.77873 (18)	0.0459 (6)
H1	0.2231	0.0031	0.7550	0.055*
C2	0.3686 (3)	0.1270 (2)	0.86777 (19)	0.0524 (6)
H2	0.3708	0.1545	0.9230	0.063*
C3	0.5430 (3)	0.0878 (2)	0.62302 (18)	0.0523 (6)
H3A	0.4815	0.0482	0.5926	0.063*
H3B	0.6517	0.0410	0.6337	0.063*
C4	0.5813 (3)	0.2132 (2)	0.54720 (17)	0.0470 (6)
H4A	0.6465	0.2512	0.5770	0.056*
H4B	0.4721	0.2608	0.5398	0.056*
C5	0.6839 (3)	0.2142 (2)	0.43876 (18)	0.0459 (6)
C6	0.7487 (3)	0.3311 (2)	0.35991 (17)	0.0426 (6)
C7	0.6996 (3)	0.4402 (2)	0.3773 (2)	0.0555 (7)
H7	0.6264	0.4421	0.4413	0.067*
C8	0.7584 (4)	0.5460 (2)	0.3002 (2)	0.0638 (7)
H8	0.7225	0.6192	0.3121	0.077*
C9	0.8682 (3)	0.5450 (3)	0.2070 (2)	0.0633 (7)
H9	0.9077	0.6173	0.1558	0.076*
C10	0.9208 (3)	0.4378 (3)	0.1883 (2)	0.0645 (8)
H10	0.9964	0.4368	0.1247	0.077*
C11	0.8608 (3)	0.3315 (2)	0.26470 (19)	0.0546 (7)
H11	0.8963	0.2587	0.2520	0.066*
C12	−0.2476 (3)	0.2397 (2)	0.90146 (18)	0.0429 (6)
H12	−0.2693	0.2090	0.8524	0.051*
C13	−0.1537 (3)	0.2650 (2)	1.02034 (19)	0.0506 (6)
H13	−0.0918	0.2519	1.0729	0.061*
C14	−0.4552 (3)	0.4253 (2)	0.8686 (2)	0.0570 (7)
H14A	−0.5458	0.4428	0.9245	0.068*
H14B	−0.5080	0.3868	0.8329	0.068*
C15	−0.3844 (3)	0.5427 (2)	0.78923 (19)	0.0509 (6)
H15A	−0.4804	0.6033	0.7789	0.061*
H15B	−0.3055	0.5701	0.8188	0.061*

C16	−0.2891 (3)	0.5332 (2)	0.6827 (2)	0.0509 (6)
C17	−0.2102 (3)	0.6418 (2)	0.59709 (19)	0.0479 (6)
C18	−0.2121 (3)	0.7519 (2)	0.6130 (2)	0.0607 (7)
H18	−0.2654	0.7599	0.6789	0.073*
C19	−0.1353 (4)	0.8486 (3)	0.5314 (2)	0.0715 (8)
H19	−0.1376	0.9217	0.5427	0.086*
C20	−0.0561 (4)	0.8392 (3)	0.4342 (2)	0.0702 (8)
H20	−0.0043	0.9054	0.3797	0.084*
C21	−0.0530 (4)	0.7320 (3)	0.4172 (2)	0.0764 (9)
H21	0.0011	0.7249	0.3510	0.092*
C22	−0.1298 (3)	0.6347 (3)	0.4978 (2)	0.0637 (7)
H22	−0.1277	0.5623	0.4851	0.076*
C23	−0.2068 (3)	−0.0969 (2)	0.87534 (17)	0.0438 (6)

Atomic displacement parameters (Å^2)

	U^{11}	U^{22}	U^{33}	U^{12}	U^{13}	U^{23}
Mn1	0.0344 (3)	0.0396 (3)	0.0321 (3)	−0.0090 (2)	−0.0024 (2)	−0.0078 (2)
S1	0.0835 (5)	0.0780 (5)	0.0643 (5)	−0.0293 (4)	−0.0104 (4)	−0.0341 (4)
O1	0.0930 (14)	0.0559 (12)	0.0517 (12)	−0.0116 (10)	0.0070 (10)	−0.0209 (10)
O2	0.0934 (14)	0.0546 (12)	0.0837 (15)	−0.0020 (10)	−0.0071 (11)	−0.0356 (11)
N1	0.0423 (11)	0.0464 (12)	0.0356 (12)	−0.0131 (9)	−0.0017 (9)	−0.0090 (9)
N2	0.0519 (12)	0.0649 (14)	0.0437 (13)	−0.0276 (11)	0.0015 (10)	−0.0180 (11)
N3	0.0439 (11)	0.0465 (11)	0.0342 (11)	−0.0125 (9)	−0.0007 (9)	−0.0115 (9)
N4	0.0461 (12)	0.0391 (11)	0.0368 (11)	−0.0063 (9)	−0.0088 (9)	−0.0091 (9)
N5	0.0671 (14)	0.0482 (13)	0.0480 (13)	0.0001 (11)	−0.0184 (11)	−0.0179 (11)
N6	0.0433 (11)	0.0417 (12)	0.0448 (12)	−0.0033 (9)	−0.0111 (9)	−0.0099 (10)
N7	0.0488 (12)	0.0645 (14)	0.0515 (13)	−0.0091 (11)	−0.0147 (10)	−0.0184 (11)
C1	0.0445 (14)	0.0532 (15)	0.0390 (14)	−0.0148 (11)	−0.0065 (11)	−0.0113 (12)
C2	0.0595 (16)	0.0612 (16)	0.0376 (14)	−0.0266 (13)	0.0029 (12)	−0.0179 (12)
C3	0.0522 (15)	0.0573 (16)	0.0398 (14)	−0.0100 (12)	0.0057 (12)	−0.0152 (12)
C4	0.0445 (14)	0.0537 (15)	0.0349 (13)	−0.0037 (11)	−0.0013 (11)	−0.0098 (12)
C5	0.0429 (14)	0.0518 (15)	0.0395 (14)	−0.0015 (12)	−0.0063 (11)	−0.0134 (12)
C6	0.0405 (13)	0.0533 (15)	0.0322 (13)	−0.0052 (11)	−0.0080 (11)	−0.0110 (11)
C7	0.0627 (17)	0.0549 (17)	0.0412 (15)	−0.0019 (13)	−0.0043 (13)	−0.0120 (13)
C8	0.081 (2)	0.0497 (16)	0.0557 (18)	−0.0065 (14)	−0.0147 (16)	−0.0108 (14)
C9	0.0639 (18)	0.0649 (19)	0.0529 (18)	−0.0242 (15)	−0.0152 (14)	0.0000 (15)
C10	0.0645 (18)	0.082 (2)	0.0379 (15)	−0.0223 (16)	0.0035 (13)	−0.0123 (15)
C11	0.0609 (16)	0.0605 (17)	0.0402 (15)	−0.0095 (13)	−0.0027 (12)	−0.0174 (13)
C12	0.0432 (14)	0.0435 (14)	0.0420 (14)	−0.0090 (11)	−0.0056 (11)	−0.0143 (12)
C13	0.0664 (17)	0.0460 (15)	0.0431 (15)	−0.0026 (13)	−0.0194 (13)	−0.0142 (12)
C14	0.0443 (15)	0.0577 (17)	0.0664 (18)	0.0029 (13)	−0.0162 (13)	−0.0167 (14)
C15	0.0532 (15)	0.0448 (14)	0.0529 (16)	0.0092 (12)	−0.0183 (13)	−0.0141 (13)
C16	0.0512 (15)	0.0448 (15)	0.0614 (17)	0.0087 (12)	−0.0222 (13)	−0.0208 (14)
C17	0.0501 (14)	0.0457 (15)	0.0496 (15)	0.0125 (12)	−0.0206 (12)	−0.0172 (12)
C18	0.0799 (19)	0.0510 (16)	0.0476 (16)	0.0053 (14)	−0.0106 (14)	−0.0177 (14)
C19	0.090 (2)	0.0491 (17)	0.069 (2)	0.0007 (15)	−0.0149 (18)	−0.0154 (16)
C20	0.0660 (18)	0.068 (2)	0.0579 (19)	0.0036 (15)	−0.0109 (15)	−0.0029 (16)

C21	0.071 (2)	0.088 (2)	0.0564 (19)	0.0139 (18)	−0.0056 (16)	−0.0210 (18)
C22	0.0677 (18)	0.0635 (18)	0.0621 (19)	0.0128 (15)	−0.0150 (15)	−0.0294 (16)
C23	0.0484 (15)	0.0493 (15)	0.0325 (13)	0.0019 (12)	−0.0090 (11)	−0.0139 (11)

Geometric parameters (Å, °)

Mn1—N7	2.207 (2)	C7—C8	1.375 (3)
Mn1—N1	2.2452 (17)	C7—H7	0.9300
Mn1—N4	2.2796 (18)	C8—C9	1.358 (4)
S1—C23	1.619 (3)	C8—H8	0.9300
O1—C5	1.212 (3)	C9—C10	1.370 (4)
O2—C16	1.216 (3)	C9—H9	0.9300
N1—C1	1.320 (3)	C10—C11	1.378 (3)
N1—C2	1.349 (3)	C10—H10	0.9300
N2—C2	1.307 (3)	C11—H11	0.9300
N2—N3	1.352 (2)	C12—H12	0.9300
N3—C1	1.323 (3)	C13—H13	0.9300
N3—C3	1.456 (3)	C14—C15	1.510 (3)
N4—C12	1.320 (3)	C14—H14A	0.9700
N4—C13	1.352 (3)	C14—H14B	0.9700
N5—C13	1.315 (3)	C15—C16	1.504 (3)
N5—N6	1.353 (2)	C15—H15A	0.9700
N6—C12	1.326 (3)	C15—H15B	0.9700
N6—C14	1.459 (3)	C16—C17	1.490 (3)
N7—C23	1.158 (3)	C17—C22	1.380 (3)
C1—H1	0.9300	C17—C18	1.393 (3)
C2—H2	0.9300	C18—C19	1.374 (4)
C3—C4	1.503 (3)	C18—H18	0.9300
C3—H3A	0.9700	C19—C20	1.361 (4)
C3—H3B	0.9700	C19—H19	0.9300
C4—C5	1.504 (3)	C20—C21	1.367 (4)
C4—H4A	0.9700	C20—H20	0.9300
C4—H4B	0.9700	C21—C22	1.373 (4)
C5—C6	1.490 (3)	C21—H21	0.9300
C6—C7	1.380 (3)	C22—H22	0.9300
C6—C11	1.383 (3)		
N7—Mn1—N7^{i}	180.0	C8—C7—H7	119.9
N7—Mn1—N1^{i}	91.32 (7)	C6—C7—H7	119.9
N7^{i}—Mn1—N1^{i}	88.68 (7)	C9—C8—C7	120.8 (3)
N7—Mn1—N1	88.68 (7)	C9—C8—H8	119.6
N7^{i}—Mn1—N1	91.32 (7)	C7—C8—H8	119.6
N1^{i}—Mn1—N1	180.0	C8—C9—C10	120.1 (3)
N7—Mn1—N4^{i}	89.00 (7)	C8—C9—H9	119.9
N7^{i}—Mn1—N4^{i}	91.00 (7)	C10—C9—H9	119.9
N1^{i}—Mn1—N4^{i}	93.23 (6)	C9—C10—C11	119.4 (2)
N1—Mn1—N4^{i}	86.77 (6)	C9—C10—H10	120.3
N7—Mn1—N4	91.00 (7)	C11—C10—H10	120.3

$N7^{i}$—Mn1—N4	89.00 (7)	C10—C11—C6	121.2 (2)
$N1^{i}$—Mn1—N4	86.77 (6)	C10—C11—H11	119.4
N1—Mn1—N4	93.23 (6)	C6—C11—H11	119.4
$N4^{i}$—Mn1—N4	180.000 (1)	N4—C12—N6	110.5 (2)
C1—N1—C2	102.08 (18)	N4—C12—H12	124.7
C1—N1—Mn1	127.56 (15)	N6—C12—H12	124.7
C2—N1—Mn1	130.23 (15)	N5—C13—N4	115.0 (2)
C2—N2—N3	102.28 (18)	N5—C13—H13	122.5
C1—N3—N2	109.70 (18)	N4—C13—H13	122.5
C1—N3—C3	129.7 (2)	N6—C14—C15	112.91 (19)
N2—N3—C3	120.58 (18)	N6—C14—H14A	109.0
C12—N4—C13	102.41 (19)	C15—C14—H14A	109.0
C12—N4—Mn1	129.36 (15)	N6—C14—H14B	109.0
C13—N4—Mn1	127.34 (15)	C15—C14—H14B	109.0
C13—N5—N6	102.19 (19)	H14A—C14—H14B	107.8
C12—N6—N5	109.89 (19)	C16—C15—C14	113.7 (2)
C12—N6—C14	130.0 (2)	C16—C15—H15A	108.8
N5—N6—C14	120.02 (19)	C14—C15—H15A	108.8
C23—N7—Mn1	143.02 (18)	C16—C15—H15B	108.8
N1—C1—N3	110.6 (2)	C14—C15—H15B	108.8
N1—C1—H1	124.7	H15A—C15—H15B	107.7
N3—C1—H1	124.7	O2—C16—C17	120.4 (2)
N2—C2—N1	115.3 (2)	O2—C16—C15	120.1 (2)
N2—C2—H2	122.3	C17—C16—C15	119.5 (2)
N1—C2—H2	122.3	C22—C17—C18	117.7 (2)
N3—C3—C4	110.74 (19)	C22—C17—C16	119.6 (2)
N3—C3—H3A	109.5	C18—C17—C16	122.7 (2)
C4—C3—H3A	109.5	C19—C18—C17	120.1 (3)
N3—C3—H3B	109.5	C19—C18—H18	119.9
C4—C3—H3B	109.5	C17—C18—H18	119.9
H3A—C3—H3B	108.1	C20—C19—C18	121.1 (3)
C3—C4—C5	112.8 (2)	C20—C19—H19	119.4
C3—C4—H4A	109.0	C18—C19—H19	119.4
C5—C4—H4A	109.0	C19—C20—C21	119.6 (3)
C3—C4—H4B	109.0	C19—C20—H20	120.2
C5—C4—H4B	109.0	C21—C20—H20	120.2
H4A—C4—H4B	107.8	C20—C21—C22	120.0 (3)
O1—C5—C6	121.1 (2)	C20—C21—H21	120.0
O1—C5—C4	120.6 (2)	C22—C21—H21	120.0
C6—C5—C4	118.3 (2)	C21—C22—C17	121.5 (3)
C7—C6—C11	118.2 (2)	C21—C22—H22	119.3
C7—C6—C5	122.6 (2)	C17—C22—H22	119.3
C11—C6—C5	119.2 (2)	N7—C23—S1	178.0 (2)
C8—C7—C6	120.3 (2)		
N7—Mn1—N1—C1	−17.61 (19)	O1—C5—C6—C7	170.4 (2)
$N7^{i}$—Mn1—N1—C1	162.39 (19)	C4—C5—C6—C7	−8.7 (3)
$N4^{i}$—Mn1—N1—C1	71.47 (19)	O1—C5—C6—C11	−8.7 (3)
N4—Mn1—N1—C1	−108.53 (19)	C4—C5—C6—C11	172.2 (2)

N7—Mn1—N1—C2	167.2 (2)	C11—C6—C7—C8	1.4 (4)
N7^{i}—Mn1—N1—C2	−12.8 (2)	C5—C6—C7—C8	−177.7 (2)
N4^{i}—Mn1—N1—C2	−103.7 (2)	C6—C7—C8—C9	−1.4 (4)
N4—Mn1—N1—C2	76.3 (2)	C7—C8—C9—C10	0.5 (4)
C2—N2—N3—C1	−0.3 (3)	C8—C9—C10—C11	0.3 (4)
C2—N2—N3—C3	−177.8 (2)	C9—C10—C11—C6	−0.1 (4)
N7—Mn1—N4—C12	2.65 (19)	C7—C6—C11—C10	−0.7 (4)
N7^{i}—Mn1—N4—C12	−177.35 (19)	C5—C6—C11—C10	178.5 (2)
N1^{i}—Mn1—N4—C12	−88.62 (19)	C13—N4—C12—N6	0.3 (2)
N1—Mn1—N4—C12	91.38 (19)	Mn1—N4—C12—N6	170.01 (13)
N7—Mn1—N4—C13	169.91 (18)	N5—N6—C12—N4	−0.2 (2)
N7^{i}—Mn1—N4—C13	−10.09 (18)	C14—N6—C12—N4	−176.4 (2)
N1^{i}—Mn1—N4—C13	78.64 (18)	N6—N5—C13—N4	0.3 (3)
N1—Mn1—N4—C13	−101.36 (18)	C12—N4—C13—N5	−0.4 (3)
C13—N5—N6—C12	0.0 (2)	Mn1—N4—C13—N5	−170.35 (15)
C13—N5—N6—C14	176.6 (2)	C12—N6—C14—C15	−107.7 (3)
N1^{i}—Mn1—N7—C23	−27.0 (3)	N5—N6—C14—C15	76.5 (3)
N1—Mn1—N7—C23	153.0 (3)	N6—C14—C15—C16	77.9 (3)
N4^{i}—Mn1—N7—C23	66.2 (3)	C14—C15—C16—O2	2.3 (3)
N4—Mn1—N7—C23	−113.8 (3)	C14—C15—C16—C17	−179.2 (2)
C2—N1—C1—N3	−0.3 (3)	O2—C16—C17—C22	1.5 (4)
Mn1—N1—C1—N3	−176.47 (14)	C15—C16—C17—C22	−177.0 (2)
N2—N3—C1—N1	0.4 (3)	O2—C16—C17—C18	−177.9 (2)
C3—N3—C1—N1	177.6 (2)	C15—C16—C17—C18	3.6 (3)
N3—N2—C2—N1	0.2 (3)	C22—C17—C18—C19	−0.2 (4)
C1—N1—C2—N2	0.0 (3)	C16—C17—C18—C19	179.2 (2)
Mn1—N1—C2—N2	176.10 (16)	C17—C18—C19—C20	−0.2 (4)
C1—N3—C3—C4	125.4 (2)	C18—C19—C20—C21	0.2 (4)
N2—N3—C3—C4	−57.7 (3)	C19—C20—C21—C22	0.1 (4)
N3—C3—C4—C5	−177.45 (19)	C20—C21—C22—C17	−0.5 (4)
C3—C4—C5—O1	8.0 (3)	C18—C17—C22—C21	0.5 (4)
C3—C4—C5—C6	−172.95 (19)	C16—C17—C22—C21	−178.9 (2)

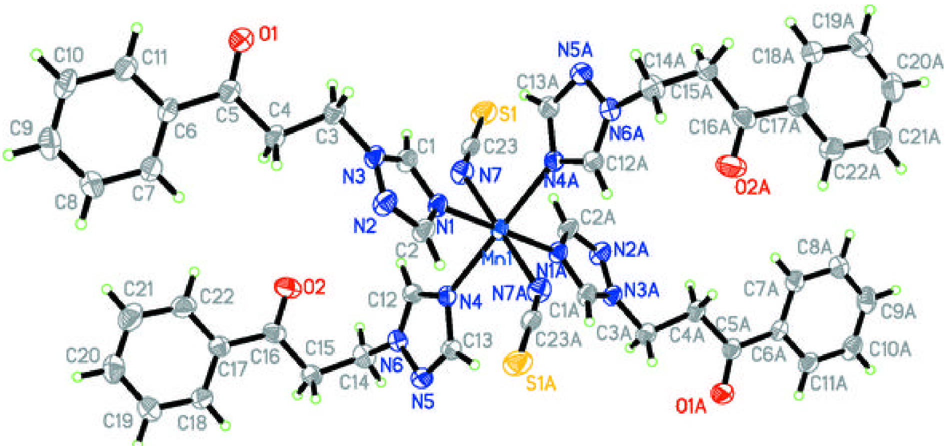
Symmetry codes: (i) −*x*, −*y*, −*z*+2.

Hydrogen-bond geometry (Å, °)

D—H···*A*	*D*—H	H···*A*	*D*···*A*	*D*—H···*A*
C12—H12···N2^{ii}	0.93	2.62	3.436 (3)	146
C18—H18···S1^{iii}	0.93	2.82	3.725 (3)	164

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

Fig. 1


O1
C10
C11
C5
C9
C6
C4
C3
C8
C7
C1
N3
N2
N1
C2
Mn1
C23
S1
N7
N5A
C13A
N6A
N4A
C12A
C14A
C15A
C16A
C17A
C18A
C19A
C20A
C21A
C22A
O2A
C2A
N2A
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N7A
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C3A
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C5A
C6A
C7A
C8A
C9A
C10A
C11A
O1A
C23A
S1A
O2
C21
C22
C20
C17
C16
C15
C14
N6
N5
C12
N4
C13
C19
C18

Fig. 2