



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 04:47 PM UTC

PDB ID : 2VZW / pdb_00002vzw
Title : X-ray structure of the heme-bound GAF domain of sensory histidine kinase
DosT of Mycobacterium tuberculosis
Authors : Podust, L.M.; Ioanoviciu, A.; Ortiz de Montellano, P.R.
Deposited on : 2008-08-06
Resolution : 2.30 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 2.0
EDS : 3.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

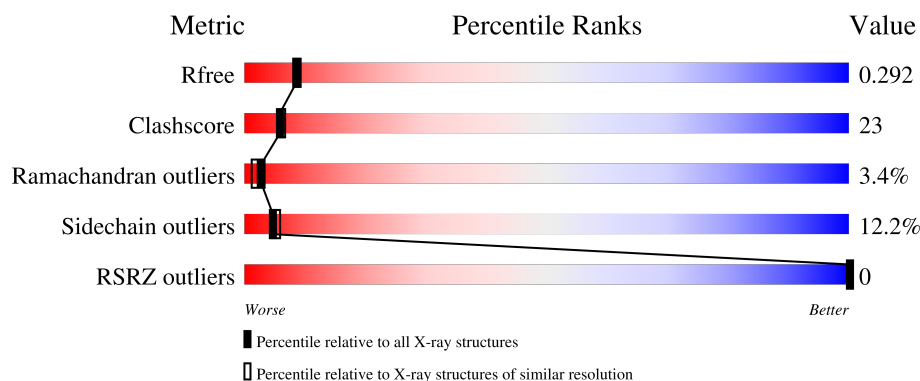
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.30 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	6319 (2.30-2.30)
Clashscore	190562	6919 (2.30-2.30)
Ramachandran outliers	187476	6854 (2.30-2.30)
Sidechain outliers	187428	6854 (2.30-2.30)
RSRZ outliers	180081	6325 (2.30-2.30)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	149	
1	B	149	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	ACT	B	1211	-	-	X	-

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 2509 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

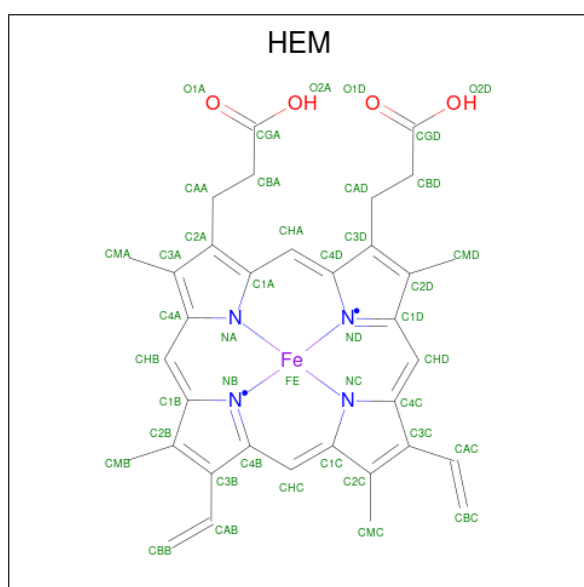
- Molecule 1 is a protein called PROBABLE HISTIDINE KINASE RESPONSE REGULATOR.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	146	Total	C	N	O	S	0	3	0
			1143	724	207	211	1			
1	B	149	Total	C	N	O	S	0	2	0
			1161	733	208	219	1			

There are 2 discrepancies between the modelled and reference sequences:

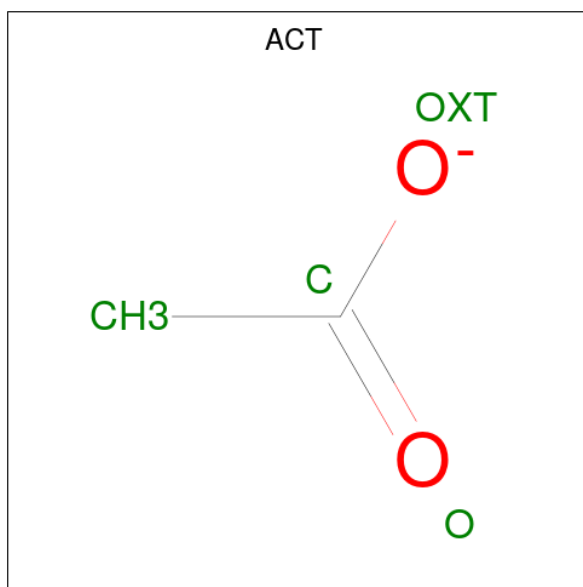
Chain	Residue	Modelled	Actual	Comment	Reference
A	60	GLY	-	expression tag	UNP O53473
B	60	GLY	-	expression tag	UNP O53473

- Molecule 2 is PROTOPORPHYRIN IX CONTAINING FE (CCD ID: HEM) (formula: $C_{34}H_{32}FeN_4O_4$).



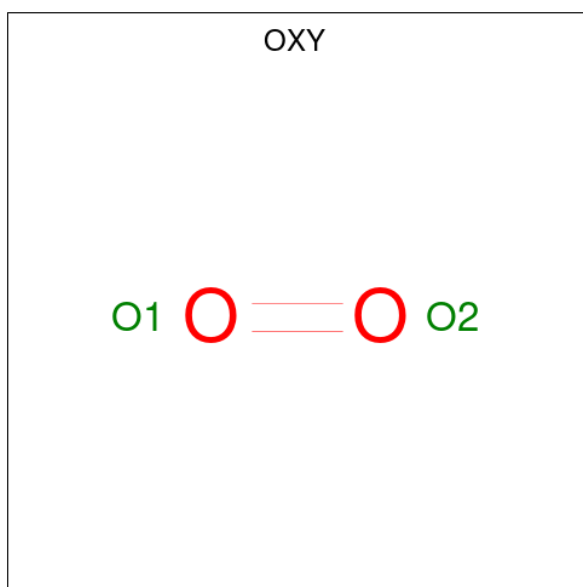
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total 43	C 34	Fe 1	N 4	O 4	0	0
2	B	1	Total 43	C 34	Fe 1	N 4	O 4	0	0

- Molecule 3 is ACETATE ION (CCD ID: ACT) (formula: $C_2H_3O_2$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			4	2	2		
3	B	1	Total	C	O	0	0
			4	2	2		

- Molecule 4 is OXYGEN MOLECULE (CCD ID: OXY) (formula: O_2).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	B	1	Total O 2 2	0	0

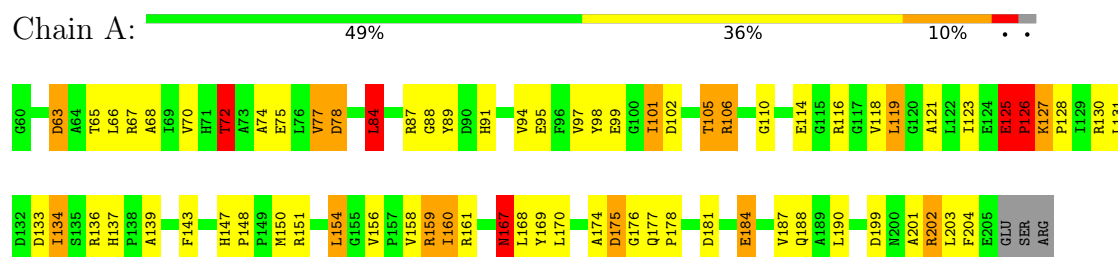
- Molecule 5 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	52	Total O 52 52	0	0
5	B	57	Total O 57 57	0	0

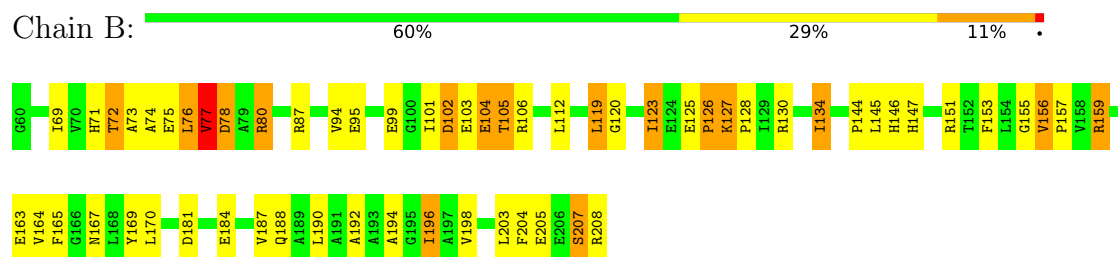
3 Residue-property plots

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

• Molecule 1: PROBABLE HISTIDINE KINASE RESPONSE REGULATOR



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4 Data and refinement statistics

Property	Value	Source
Space group	P 43	Depositor
Cell constants a, b, c, α , β , γ	87.92Å 87.92Å 66.76Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	88.05 – 2.30 87.92 – 2.30	Depositor EDS
% Data completeness (in resolution range)	100.0 (88.05-2.30) 99.9 (87.92-2.30)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.26 (at 2.29Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.236 , 0.290 0.243 , 0.292	Depositor DCC
R_{free} test set	1171 reflections (5.11%)	wwPDB-VP
Wilson B-factor (Å ²)	47.2	Xtriage
Anisotropy	0.407	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 38.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.40$, $\langle L^2 \rangle = 0.23$	Xtriage
Estimated twinning fraction	0.407 for h,-k,-l	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	2509	wwPDB-VP
Average B, all atoms (Å ²)	49.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.61% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: ACT, HEM, OXY

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z > 5$	RMSZ	$\# Z > 5$
1	A	1.18	0/1167	1.60	19/1588 (1.2%)
1	B	1.30	3/1185 (0.3%)	1.45	12/1611 (0.7%)
All	All	1.24	3/2352 (0.1%)	1.53	31/3199 (1.0%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	123	ILE	CA-CB	7.98	1.64	1.54
1	B	72	THR	CA-CB	5.92	1.62	1.53
1	B	77	VAL	CA-C	5.54	1.60	1.52

All (31) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	125	GLU	CA-C-N	18.23	142.63	119.84
1	A	125	GLU	C-N-CA	18.23	142.63	119.84
1	A	77	VAL	N-CA-C	-10.38	103.03	113.10
1	A	147	HIS	CA-C-N	-8.68	111.44	120.38
1	A	147	HIS	C-N-CA	-8.68	111.44	120.38
1	B	104	GLU	N-CA-C	-7.78	99.77	110.35
1	B	101	ILE	N-CA-C	7.29	119.33	108.46
1	B	119	LEU	CA-CB-CG	-6.88	92.20	116.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	147	HIS	N-CA-C	-6.83	101.46	109.93
1	B	78	ASP	N-CA-C	6.78	121.13	112.86
1	A	159	ARG	N-CA-C	6.62	119.39	108.99
1	A	94	VAL	N-CA-C	-6.50	106.22	111.81
1	B	94[A]	VAL	N-CA-C	-6.45	106.26	111.81
1	B	94[B]	VAL	N-CA-C	-6.45	106.26	111.81
1	A	199	ASP	N-CA-C	-6.31	104.33	111.14
1	B	112	LEU	CA-C-N	-6.16	113.43	119.78
1	B	112	LEU	C-N-CA	-6.16	113.43	119.78
1	B	165	PHE	N-CA-C	-6.12	105.94	113.41
1	A	77	VAL	CB-CA-C	5.93	117.46	110.93
1	A	74	ALA	N-CA-C	-5.90	104.77	111.14
1	A	167	ASN	N-CA-C	5.76	117.82	109.07
1	A	84	LEU	CA-CB-CG	5.42	135.27	116.30
1	A	78	ASP	N-CA-C	5.27	118.66	111.39
1	A	127	LYS	N-CA-C	-5.23	103.10	110.31
1	A	72	THR	N-CA-C	5.21	116.65	111.07
1	A	184	GLU	N-CA-C	5.14	116.89	111.28
1	B	196	ILE	CB-CA-C	-5.10	105.44	111.97
1	A	126	PRO	CA-N-CD	-5.09	104.88	112.00
1	B	159	ARG	N-CA-C	5.08	116.56	108.79
1	A	143	PHE	CA-C-N	5.03	125.24	120.31
1	A	143	PHE	C-N-CA	5.03	125.24	120.31

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	125	GLU	Peptide

5.2 Too-close contacts ⓘ

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1143	0	1124	61	0
1	B	1161	0	1135	47	0
2	A	43	0	30	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	B	43	0	30	3	0
3	A	4	0	3	1	0
3	B	4	0	3	2	0
4	B	2	0	0	0	0
5	A	52	0	0	3	0
5	B	57	0	0	7	0
All	All	2509	0	2325	109	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 23.

All (109) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:133:ASP:O	1:A:136[B]:ARG:HG2	1.57	1.04
1:A:66:LEU:HD12	1:A:95:GLU:HG2	1.42	0.98
1:A:98:TYR:OH	1:A:106:ARG:HG3	1.66	0.95
1:A:98:TYR:CZ	1:A:106:ARG:HG3	2.09	0.88
1:B:125:GLU:HB2	5:B:2025:HOH:O	1.73	0.87
1:A:116:ARG:NH2	5:A:2024:HOH:O	2.05	0.82
1:A:66:LEU:CD1	1:A:95:GLU:HG2	2.11	0.81
1:A:175:ASP:O	1:A:177[A]:GLN:N	2.20	0.75
1:A:154:LEU:HD13	1:A:187:VAL:HG21	1.70	0.74
1:B:192:ALA:O	1:B:196:ILE:HG12	1.89	0.72
1:B:125:GLU:O	1:B:126:PRO:O	2.08	0.72
1:A:106:ARG:HE	1:A:106:ARG:C	1.97	0.71
1:A:63:ASP:O	1:A:67:ARG:HG3	1.92	0.69
1:A:119:LEU:HD21	1:A:169:TYR:CZ	2.28	0.69
1:A:75:GLU:HG2	1:B:76:LEU:HD13	1.74	0.68
1:B:203:LEU:O	5:B:2056:HOH:O	2.12	0.68
1:A:72:THR:HG22	1:A:190:LEU:CD1	2.24	0.67
1:A:137:HIS:HD2	1:A:139:ALA:H	1.42	0.67
1:A:67:ARG:HA	1:A:97[B]:VAL:HG11	1.76	0.66
1:A:175:ASP:O	1:A:177[B]:GLN:N	2.26	0.66
1:B:120:GLY:HA2	1:B:123:ILE:HD12	1.77	0.66
1:B:95:GLU:OE1	5:B:2012:HOH:O	2.13	0.66
1:A:175:ASP:C	1:A:177[A]:GLN:H	2.05	0.65
1:A:89:TYR:OH	1:A:204:PHE:HA	1.96	0.65
1:A:87:ARG:HH12	3:A:1207:ACT:H1	1.62	0.64
1:B:128:PRO:HB3	1:B:156:VAL:HG12	1.81	0.63
1:B:157:PRO:HB3	5:B:2024:HOH:O	1.99	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:87:ARG:HH12	3:B:1211:ACT:H2	1.63	0.61
1:B:127:LYS:CB	5:B:2026:HOH:O	2.48	0.61
1:B:159:ARG:HH11	1:B:164:VAL:CG2	2.13	0.61
1:A:160:ILE:HG22	1:A:161:ARG:HG3	1.82	0.61
1:B:87:ARG:HH12	3:B:1211:ACT:CH3	2.14	0.61
1:B:159:ARG:HH11	1:B:164:VAL:HG22	1.66	0.61
2:B:1209:HEM:HHC	2:B:1209:HEM:HBB2	1.83	0.60
1:A:119:LEU:HD21	1:A:169:TYR:CE2	2.37	0.59
1:A:72:THR:HG22	1:A:190:LEU:HD13	1.82	0.59
1:A:88:GLY:HA2	1:A:204:PHE:HE1	1.68	0.59
1:B:145:LEU:O	1:B:146:HIS:HB2	2.03	0.59
1:A:175:ASP:C	1:A:177[B]:GLN:H	2.10	0.59
1:B:123:ILE:HD11	5:B:2045:HOH:O	2.03	0.59
2:B:1209:HEM:HMC1	2:B:1209:HEM:HBC2	1.85	0.59
1:B:74:ALA:HA	1:B:170:LEU:HD22	1.84	0.58
1:B:72:THR:O	1:B:76:LEU:HB2	2.04	0.57
1:B:159:ARG:HB2	1:B:163:GLU:O	2.05	0.57
1:A:169:TYR:C	1:A:170:LEU:HD12	2.30	0.57
2:A:1206:HEM:HBB2	2:A:1206:HEM:HHC	1.86	0.56
1:B:126:PRO:O	1:B:127:LYS:O	2.24	0.56
1:A:101:ILE:HD11	1:A:148:PRO:HG2	1.88	0.55
1:A:102:ASP:OD1	1:A:105:THR:HG23	2.07	0.55
1:B:120:GLY:HA2	1:B:123:ILE:HB	1.89	0.55
1:B:204:PHE:O	1:B:207:SER:HB2	2.07	0.55
1:A:134:ILE:HD11	1:A:150:MET:O	2.08	0.54
1:A:184:GLU:O	1:A:188:GLN:HG3	2.07	0.54
1:A:72:THR:HG22	1:A:190:LEU:HD11	1.90	0.54
1:A:101:ILE:HD11	1:A:148:PRO:CG	2.37	0.54
1:A:70:VAL:CG2	1:A:84:LEU:HB3	2.37	0.54
1:A:72:THR:HG23	1:B:190:LEU:HD21	1.91	0.52
1:A:106:ARG:NH2	1:A:110:GLY:O	2.43	0.52
1:A:125:GLU:C	1:A:127:LYS:H	2.18	0.52
1:A:174:ALA:O	1:A:175:ASP:O	2.28	0.52
1:B:106:ARG:HD3	1:B:106:ARG:C	2.35	0.51
1:B:104:GLU:O	1:B:105:THR:CB	2.59	0.51
1:B:130:ARG:NH2	1:B:181[A]:ASP:OD1	2.42	0.51
1:A:68:ALA:O	1:A:72:THR:HB	2.12	0.49
1:A:84:LEU:HD13	1:A:158:VAL:HG21	1.95	0.49
1:A:127:LYS:HG2	1:A:128:PRO:O	2.13	0.48
1:B:144:PRO:HG3	2:B:1209:HEM:HMD3	1.96	0.48
1:A:84:LEU:HA	1:A:167:ASN:O	2.14	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:167:ASN:ND2	5:B:2045:HOH:O	2.46	0.47
1:B:74:ALA:HA	1:B:170:LEU:CD2	2.44	0.47
1:A:88:GLY:HA2	1:A:204:PHE:CE1	2.50	0.47
1:A:131:LEU:HD13	1:A:134:ILE:HA	1.96	0.47
1:B:77:VAL:HG21	1:B:187:VAL:HG23	1.97	0.46
1:B:134:ILE:HG13	1:B:153:PHE:HB2	1.98	0.46
1:A:134:ILE:HG13	1:A:151:ARG:O	2.15	0.45
1:B:159:ARG:NH1	1:B:164:VAL:CG2	2.79	0.45
1:A:177[A]:GLN:HG3	1:A:178:PRO:HD2	1.99	0.45
1:A:201:ALA:O	1:A:202:ARG:HB2	2.16	0.45
1:B:205:GLU:OE2	1:B:208:ARG:NH2	2.45	0.45
1:B:102:ASP:OD1	1:B:105:THR:OG1	2.26	0.44
1:A:70:VAL:HG23	1:A:84:LEU:HD23	1.99	0.44
1:B:103:GLU:C	1:B:104:GLU:O	2.59	0.44
1:B:159:ARG:NH1	1:B:164:VAL:HG23	2.32	0.44
1:A:151:ARG:HE	1:A:151:ARG:HB2	1.45	0.44
1:A:160:ILE:HG22	1:A:161:ARG:N	2.33	0.44
1:B:80:ARG:CG	1:B:80:ARG:HH11	2.31	0.43
1:B:184:GLU:O	1:B:188:GLN:HG3	2.18	0.43
1:A:118:VAL:O	1:A:121:ALA:HB3	2.18	0.43
1:B:71:HIS:ND1	1:B:99:GLU:OE1	2.34	0.43
1:B:145:LEU:O	1:B:146:HIS:CB	2.67	0.43
1:A:134:ILE:HG13	1:A:134:ILE:H	1.48	0.43
1:A:160:ILE:CG2	1:A:161:ARG:N	2.81	0.43
1:A:67:ARG:HB3	1:A:67:ARG:NH1	2.34	0.42
1:A:70:VAL:HG21	1:A:84:LEU:HB3	2.01	0.42
1:B:194:ALA:O	1:B:198:VAL:HG23	2.19	0.42
1:A:105:THR:CG2	5:A:2013:HOH:O	2.66	0.42
1:A:87:ARG:HD3	1:A:91:HIS:O	2.19	0.42
1:A:128:PRO:HB3	1:A:156:VAL:HG12	2.02	0.42
1:B:69:ILE:HG23	1:B:190:LEU:HB3	2.02	0.42
1:B:128:PRO:HA	1:B:155:GLY:O	2.19	0.42
1:A:114:GLU:OE1	1:A:116:ARG:NH2	2.49	0.41
1:B:104:GLU:O	1:B:105:THR:HB	2.20	0.41
1:B:127:LYS:HA	1:B:128:PRO:HD3	1.86	0.41
1:A:130:ARG:NH2	1:A:181:ASP:OD1	2.53	0.41
1:B:169:TYR:C	1:B:170:LEU:HG	2.46	0.41
1:A:105:THR:HG23	5:A:2013:HOH:O	2.19	0.41
1:A:174:ALA:C	1:A:175:ASP:O	2.64	0.41
1:A:98:TYR:HH	1:A:106:ARG:HG3	1.77	0.41
1:B:73:ALA:O	1:B:77:VAL:HB	2.22	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	147/149 (99%)	140 (95%)	2 (1%)	5 (3%)	3	2
1	B	149/149 (100%)	134 (90%)	10 (7%)	5 (3%)	3	2
All	All	296/298 (99%)	274 (93%)	12 (4%)	10 (3%)	3	2

All (10) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	126	PRO
1	A	175	ASP
1	B	126	PRO
1	B	127	LYS
1	A	176	GLY
1	A	202	ARG
1	A	203	LEU
1	B	105	THR
1	B	207	SER
1	B	102	ASP

5.3.2 Protein sidechains [i](#)

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	116/119 (98%)	97 (84%)	19 (16%)	2	2

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	B	119/119 (100%)	110 (92%)	9 (8%)	12	16
All	All	235/238 (99%)	207 (88%)	28 (12%)	5	6

All (28) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	63	ASP
1	A	65	THR
1	A	72	THR
1	A	77	VAL
1	A	78	ASP
1	A	84	LEU
1	A	99	GLU
1	A	101	ILE
1	A	105	THR
1	A	106	ARG
1	A	119	LEU
1	A	123	ILE
1	A	126	PRO
1	A	134	ILE
1	A	154	LEU
1	A	159	ARG
1	A	160	ILE
1	A	167	ASN
1	A	168	LEU
1	B	75	GLU
1	B	76	LEU
1	B	77	VAL
1	B	78	ASP
1	B	80	ARG
1	B	119	LEU
1	B	134	ILE
1	B	151	ARG
1	B	156	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	91	HIS
1	A	107	HIS
1	A	137	HIS

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Mol	Chain	Res	Type
1	A	146	HIS
1	A	167	ASN
1	A	188	GLN
1	A	200	ASN
1	B	107	HIS
1	B	146	HIS
1	B	200	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

5 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
4	OXY	B	1210	2	1,1,1	0.08	0	-		
3	ACT	A	1207	-	3,3,3	0.88	0	3,3,3	0.97	0
3	ACT	B	1211	-	3,3,3	0.64	0	3,3,3	0.87	0
2	HEM	A	1206	1	50,50,50	2.01	11 (22%)	67,82,82	1.97	23 (34%)
2	HEM	B	1209	4,1	50,50,50	1.91	9 (18%)	67,82,82	1.43	7 (10%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	HEM	A	1206	1	-	4/14/54/54	-
2	HEM	B	1209	4,1	-	1/14/54/54	-

All (20) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	1206	HEM	FE-NB	7.69	2.18	1.94
2	B	1209	HEM	C3D-C2D	7.12	1.52	1.36
2	A	1206	HEM	C3D-C2D	6.29	1.50	1.36
2	B	1209	HEM	FE-ND	5.42	2.11	1.94
2	A	1206	HEM	FE-ND	4.32	2.08	1.94
2	B	1209	HEM	FE-NC	4.04	2.08	1.95
2	A	1206	HEM	CAB-C3B	3.53	1.56	1.47
2	A	1206	HEM	CAC-C3C	2.96	1.55	1.47
2	B	1209	HEM	CAC-C3C	2.88	1.55	1.47
2	B	1209	HEM	CAB-C3B	2.85	1.55	1.47
2	B	1209	HEM	C1B-NB	-2.70	1.35	1.40
2	A	1206	HEM	CMD-C2D	2.67	1.56	1.50
2	B	1209	HEM	FE-NB	2.64	2.03	1.94
2	A	1206	HEM	FE-NA	2.55	2.03	1.95
2	B	1209	HEM	FE-NA	2.50	2.03	1.95
2	A	1206	HEM	CMB-C2B	2.49	1.55	1.50
2	B	1209	HEM	CMC-C2C	2.46	1.55	1.50
2	A	1206	HEM	C3C-C4C	-2.42	1.41	1.46
2	A	1206	HEM	CMA-C3A	2.35	1.55	1.50
2	A	1206	HEM	C2A-C3A	-2.04	1.33	1.38

All (30) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	1209	HEM	CAA-CBA-CGA	-4.80	100.94	113.67
2	B	1209	HEM	C4D-ND-C1D	4.79	110.89	105.21
2	A	1206	HEM	CHD-C4C-NC	4.09	128.90	124.45
2	A	1206	HEM	C4D-ND-C1D	3.94	109.88	105.21
2	B	1209	HEM	CMD-C2D-C1D	3.85	131.06	125.03
2	A	1206	HEM	CAD-C3D-C4D	3.85	131.40	124.70
2	A	1206	HEM	C2A-C1A-NA	-3.85	105.88	110.15
2	A	1206	HEM	C3B-C4B-NB	-3.72	106.80	109.47

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	1206	HEM	C1B-NB-C4B	3.46	109.30	105.21
2	A	1206	HEM	CAD-C3D-C2D	-3.44	121.42	127.87
2	A	1206	HEM	CAC-C3C-C4C	-3.37	116.78	124.82
2	A	1206	HEM	CAA-C2A-C1A	3.23	131.25	124.94
2	A	1206	HEM	CAA-C2A-C3A	-3.11	120.11	127.07
2	B	1209	HEM	CHA-C4D-ND	2.93	127.99	124.37
2	A	1206	HEM	O1A-CGA-CBA	-2.79	114.25	123.09
2	A	1206	HEM	CBD-CAD-C3D	-2.75	104.94	112.53
2	A	1206	HEM	C3A-C4A-NA	-2.71	105.79	110.14
2	A	1206	HEM	CHB-C4A-NA	2.71	128.77	123.86
2	A	1206	HEM	C1D-C2D-C3D	-2.69	104.15	106.98
2	B	1209	HEM	C1D-C2D-C3D	-2.68	104.16	106.98
2	A	1206	HEM	CHC-C4B-C3B	2.62	130.30	125.07
2	A	1206	HEM	C4A-NA-C1A	2.62	110.09	105.82
2	A	1206	HEM	CMD-C2D-C1D	2.59	129.09	125.03
2	A	1206	HEM	C3D-C4D-ND	-2.44	107.50	110.17
2	A	1206	HEM	C4A-C3A-C2A	2.42	109.58	106.82
2	A	1206	HEM	CAD-CBD-CGD	2.39	120.01	113.67
2	A	1206	HEM	CAC-C3C-C2C	2.20	135.60	128.43
2	B	1209	HEM	CAD-CBD-CGD	2.17	119.43	113.67
2	A	1206	HEM	C4C-NC-C1C	2.06	109.18	105.82
2	B	1209	HEM	C3D-C4D-ND	-2.04	107.93	110.17

There are no chirality outliers.

All (5) torsion outliers are listed below:

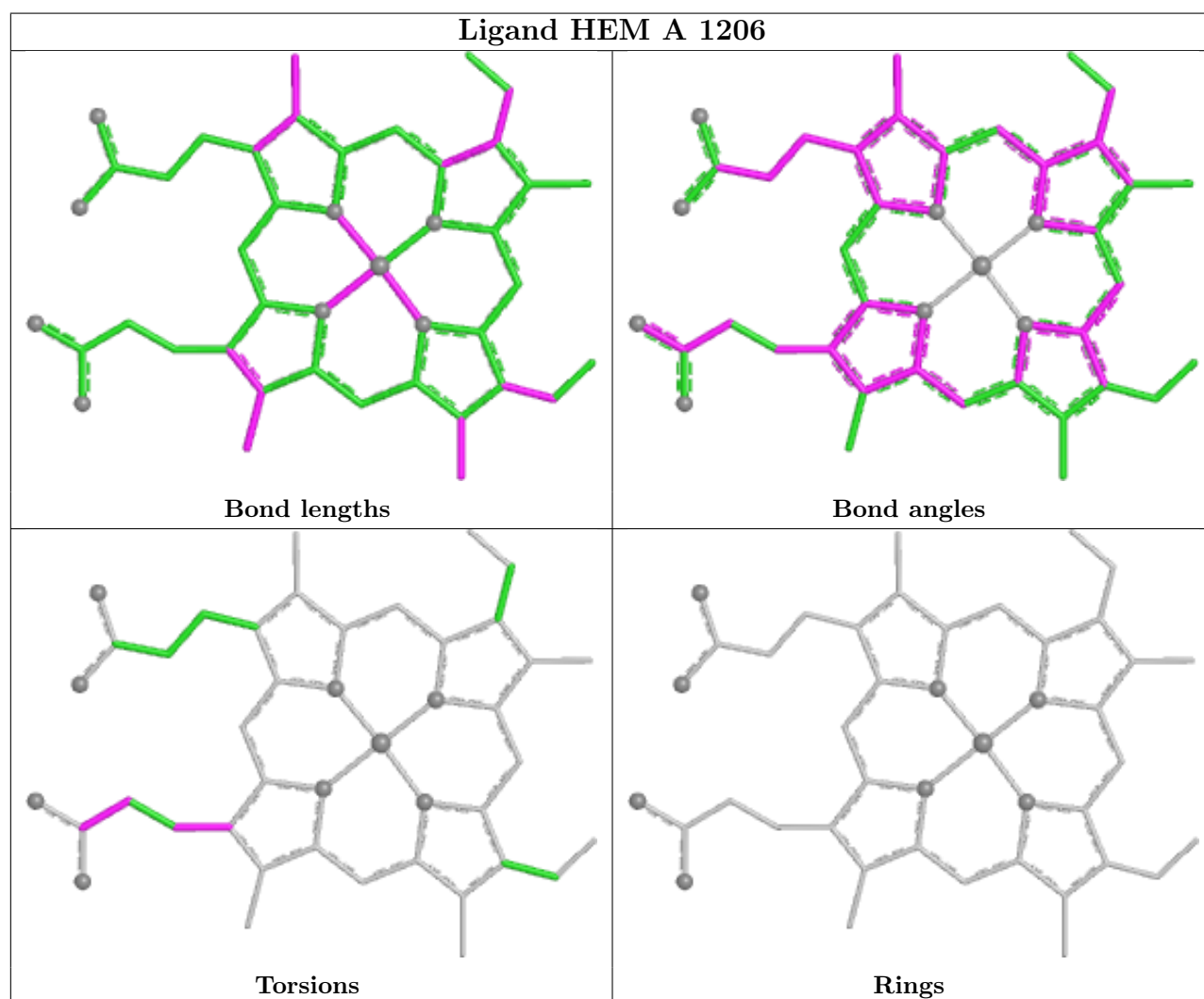
Mol	Chain	Res	Type	Atoms
2	A	1206	HEM	C1A-C2A-CAA-CBA
2	A	1206	HEM	C3A-C2A-CAA-CBA
2	A	1206	HEM	CAA-CBA-CGA-O1A
2	A	1206	HEM	CAA-CBA-CGA-O2A
2	B	1209	HEM	CAD-CBD-CGD-O2D

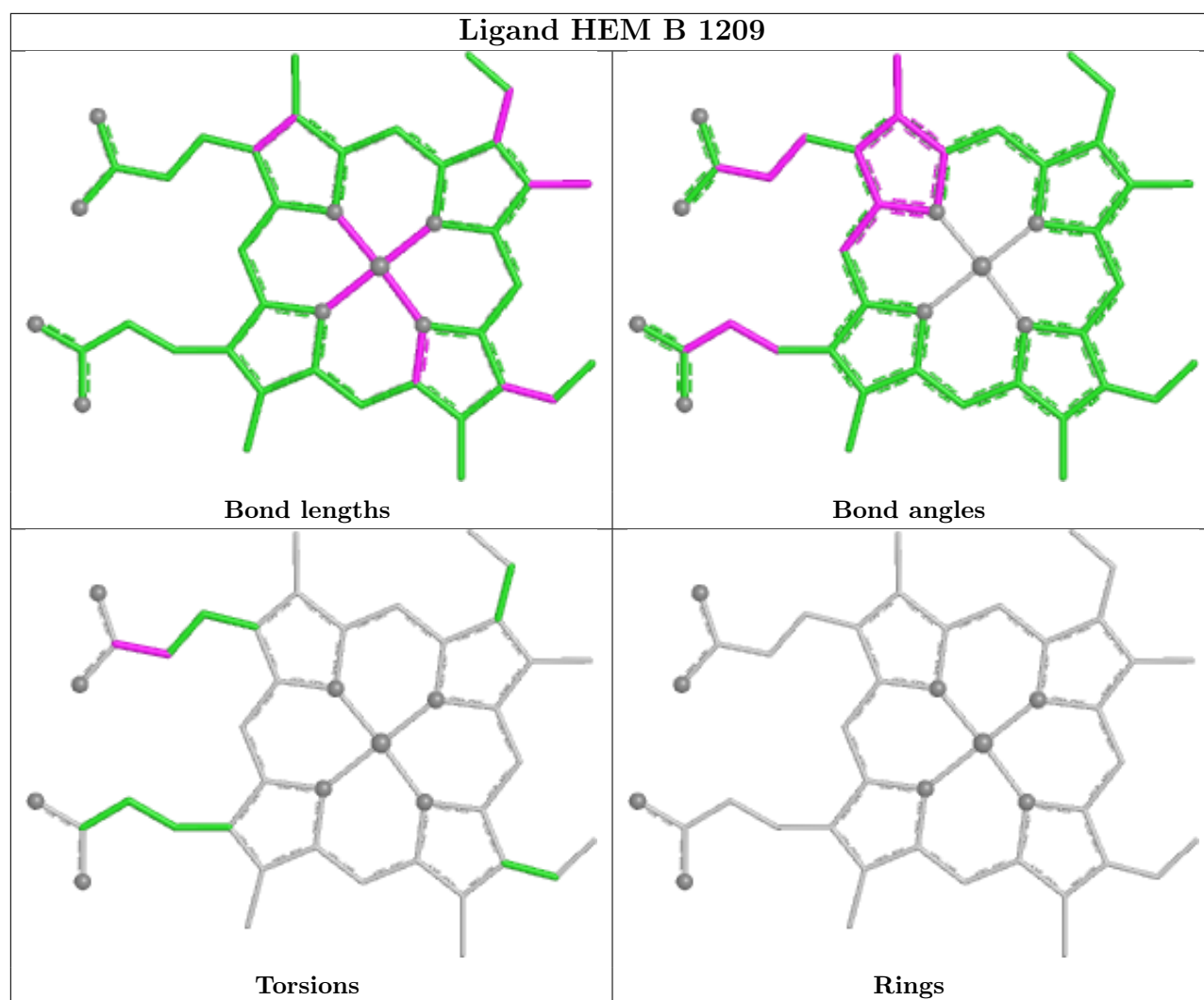
There are no ring outliers.

4 monomers are involved in 7 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	1207	ACT	1	0
3	B	1211	ACT	2	0
2	A	1206	HEM	1	0
2	B	1209	HEM	3	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.





5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data [i](#)

6.1 Protein, DNA and RNA chains [i](#)

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	146/149 (97%)	-1.12	0 100 100	23, 50, 63, 68	3 (2%)
1	B	149/149 (100%)	-1.14	0 100 100	22, 48, 63, 74	2 (1%)
All	All	295/298 (98%)	-1.13	0 100 100	22, 49, 63, 74	5 (1%)

There are no RSRZ outliers to report.

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled ‘Q< 0.9’ lists the number of atoms with occupancy less than 0.9.

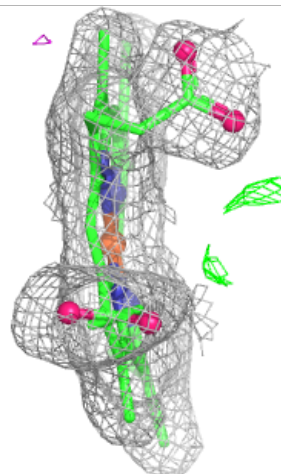
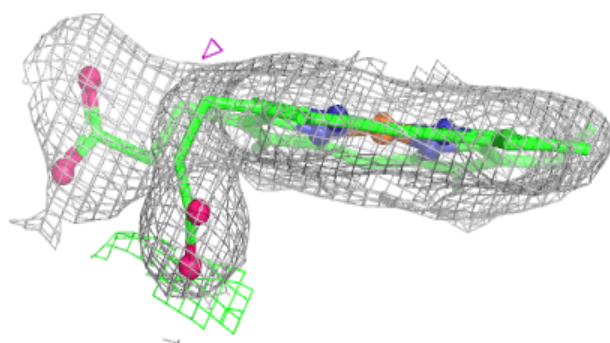
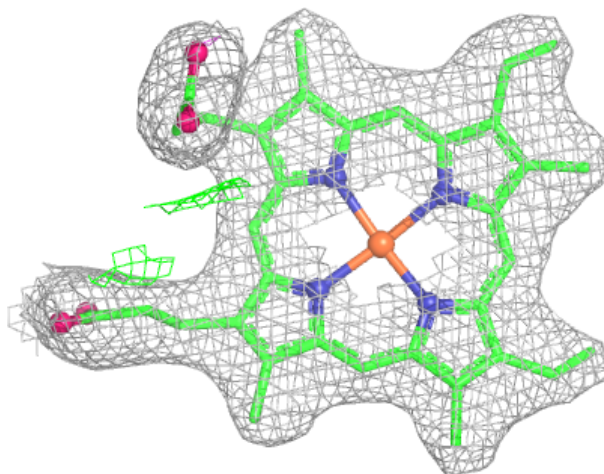
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
3	ACT	A	1207	4/4	0.98	0.07	47,51,52,53	0
3	ACT	B	1211	4/4	0.98	0.07	47,51,53,56	0
4	OXY	B	1210	2/2	0.99	0.06	42,42,42,44	0
2	HEM	B	1209	43/43	1.00	0.03	33,40,46,47	0
2	HEM	A	1206	43/43	1.00	0.04	22,33,37,39	0

The following is a graphical depiction of the model fit to experimental electron density of all

instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

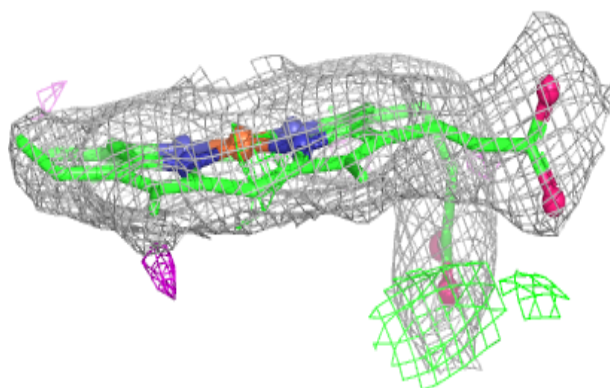
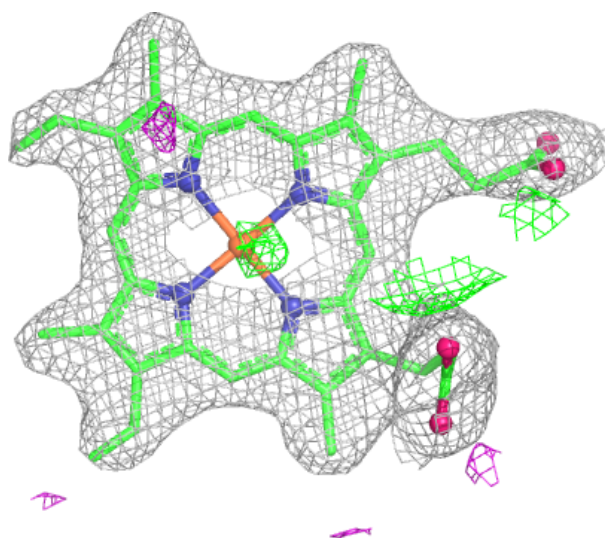
Electron density around HEM B 1209:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around HEM A 1206:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



6.5 Other polymers [i](#)

There are no such residues in this entry.