



Full wwPDB X-ray Structure Validation Report ⓘ

May 24, 2020 – 12:59 am BST

PDB ID : 1QJS
Title : mammalian blood serum haemopexin glycosylated-native protein and in complex with its ligand haem
Authors : Paoli, M.; Baker, H.M.; Morgan, W.T.; Smith, A.; Baker, E.N.
Deposited on : 1999-07-01
Resolution : 2.90 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 following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.11
buster-report : 1.1.7 (2018)
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.11

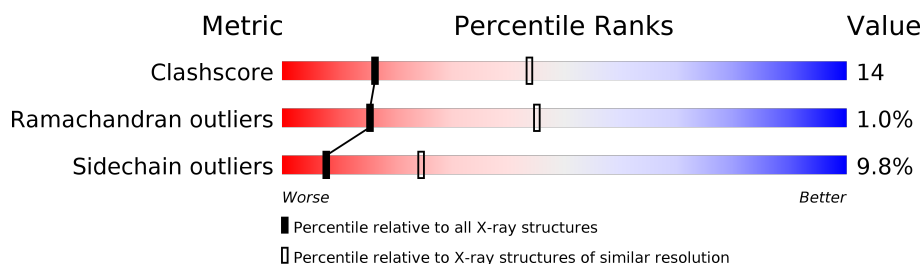
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.90 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))
Clashscore	141614	2172 (2.90-2.90)
Ramachandran outliers	138981	2115 (2.90-2.90)
Sidechain outliers	138945	2117 (2.90-2.90)

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 on 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 $\leq 5\%$.

Mol	Chain	Length	Quality of chain
1	A	460	
1	B	460	

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
2	HEM	B	500	-	-	X	-
3	PO4	B	501	-	-	X	-

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 6686 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 HEMOPEXIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	408	Total	C	N	O	S	0	0	0
			3289	2102	579	592	16			
1	B	408	Total	C	N	O	S	0	0	0
			3289	2102	579	592	16			

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



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

- Molecule 3 is PHOSPHATE ION (three-letter code: PO4) (formula: O_4P).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	O	P	0	0
			5	4	1		
3	B	1	Total	O	P	0	0
			5	4	1		

- Molecule 4 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	B	2	Total	Cl	0	0
			2	2		
4	A	2	Total	Cl	0	0
			2	2		

- Molecule 5 is SODIUM ION (three-letter code: NA) (formula: Na).

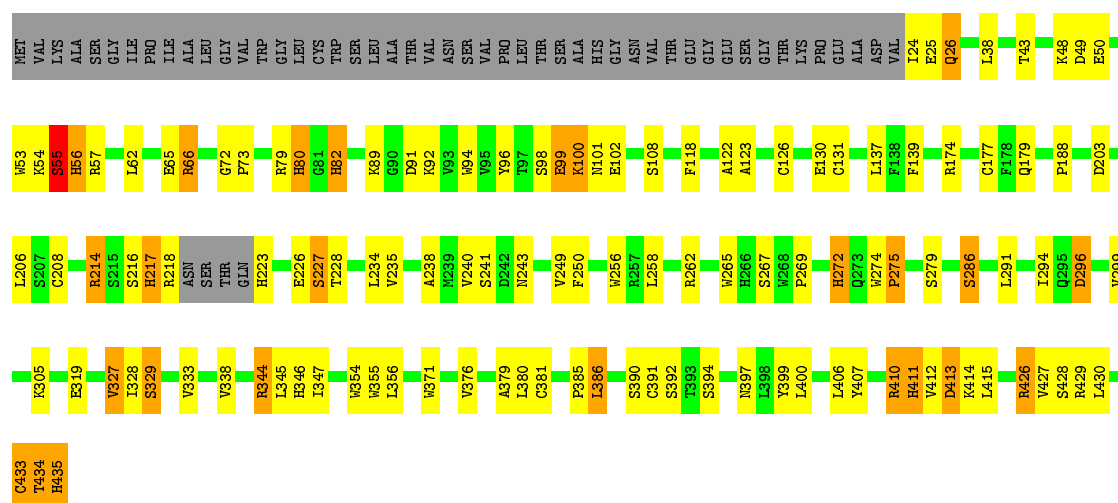
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	B	4	Total	Na	0	0
			4	4		
5	A	4	Total	Na	0	0
			4	4		

3 Residue-property plots

These plots are drawn for all protein, RNA and DNA 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. 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. 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.

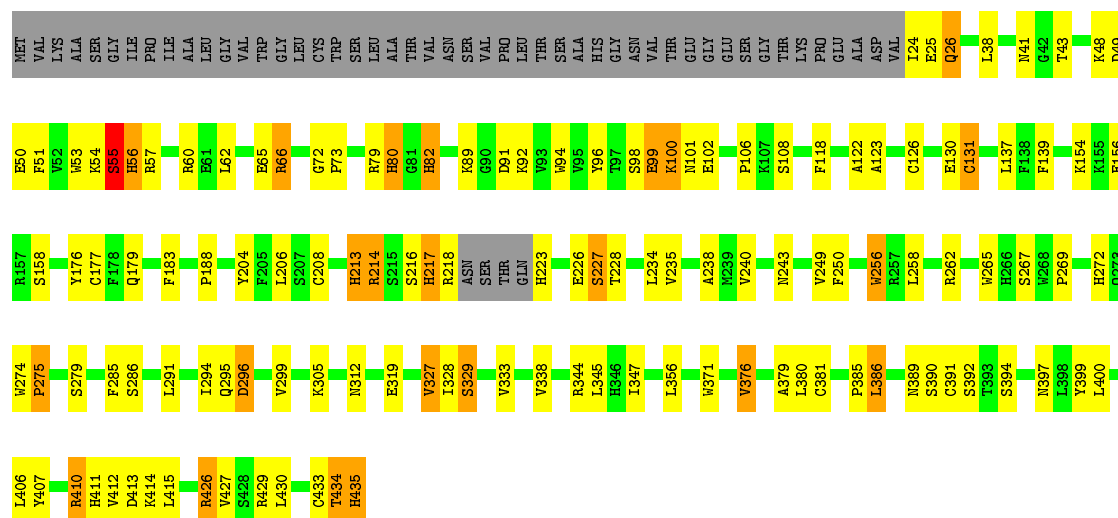
• Molecule 1: HEMOPEXIN

Chain A: 



• Molecule 1: HEMOPEXIN

Chain B: 



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	103.90 Å 69.90 Å 151.81 Å 90.00° 108.16° 90.00°	Depositor
Resolution (Å)	20.00 – 2.90 19.90 – 2.96	Depositor EDS
% Data completeness (in resolution range)	97.5 (20.00-2.90) 96.3 (19.90-2.96)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.10 (at 2.98 Å)	Xtriage
Refinement program	REFMAC	Depositor
R, R_{free}	0.255 , 0.312 0.360 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å ²)	59.9	Xtriage
Anisotropy	0.478	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.28 , 23.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.019 for h,-k,-h-l	Xtriage
F_o, F_c correlation	0.78	EDS
Total number of atoms	6686	wwPDB-VP
Average B, all atoms (Å ²)	39.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.39% 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: HEM, PO4, CL, NA

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	0.74	2/3400 (0.1%)	1.39	27/4624 (0.6%)
1	B	0.71	4/3400 (0.1%)	1.28	28/4624 (0.6%)
All	All	0.72	6/6800 (0.1%)	1.33	55/9248 (0.6%)

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	435	HIS	N-CA	-19.25	1.07	1.46
1	B	435	HIS	N-CA	-10.09	1.26	1.46
1	A	434	THR	C-N	-9.07	1.13	1.34
1	B	434	THR	CA-C	-8.59	1.30	1.52
1	B	434	THR	C-N	-6.05	1.20	1.34
1	B	433	CYS	C-N	-5.19	1.22	1.34

All (55) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	434	THR	O-C-N	29.98	170.66	122.70
1	A	434	THR	CA-C-N	-25.70	60.66	117.20
1	A	434	THR	C-N-CA	9.55	145.57	121.70
1	B	126	CYS	CA-CB-SG	-9.34	97.19	114.00
1	A	126	CYS	CA-CB-SG	-9.23	97.39	114.00
1	B	434	THR	C-N-CA	-9.06	99.06	121.70
1	A	433	CYS	C-N-CA	-8.85	99.59	121.70
1	A	386	LEU	N-CA-C	-7.67	90.29	111.00
1	B	386	LEU	N-CA-C	-7.67	90.29	111.00
1	A	381	CYS	CA-CB-SG	-7.61	100.31	114.00
1	B	208	CYS	CA-CB-SG	-7.04	101.32	114.00
1	B	381	CYS	CA-CB-SG	-7.03	101.35	114.00
1	B	433	CYS	O-C-N	-7.03	111.46	122.70

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	227	SER	CA-CB-OG	6.51	128.78	111.20
1	B	433	CYS	C-N-CA	-6.27	106.04	121.70
1	B	79	ARG	CA-C-N	6.10	130.62	117.20
1	A	275	PRO	N-CA-C	6.10	127.95	112.10
1	B	275	PRO	N-CA-C	6.05	127.84	112.10
1	B	38	LEU	O-C-N	-5.95	113.18	122.70
1	A	427	VAL	CG1-CB-CG2	-5.87	101.50	110.90
1	A	80	HIS	CA-C-N	-5.87	104.46	116.20
1	B	338	VAL	N-CA-C	-5.86	95.17	111.00
1	A	38	LEU	O-C-N	-5.81	113.41	122.70
1	B	319	GLU	CA-C-N	-5.80	104.44	117.20
1	A	319	GLU	CA-C-N	-5.78	104.49	117.20
1	B	80	HIS	CA-C-N	-5.76	104.68	116.20
1	A	79	ARG	CA-C-N	5.71	129.76	117.20
1	A	413	ASP	O-C-N	-5.70	113.57	122.70
1	A	286	SER	CA-CB-OG	5.70	126.59	111.20
1	A	391	CYS	O-C-N	-5.70	113.58	122.70
1	B	391	CYS	O-C-N	-5.68	113.62	122.70
1	A	130	GLU	N-CA-C	-5.67	95.68	111.00
1	A	296	ASP	CA-C-N	5.66	129.66	117.20
1	B	227	SER	CA-CB-OG	5.58	126.28	111.20
1	A	208	CYS	CA-CB-SG	-5.52	104.06	114.00
1	B	130	GLU	N-CA-C	-5.46	96.27	111.00
1	B	413	ASP	O-C-N	-5.43	114.02	122.70
1	B	79	ARG	O-C-N	-5.42	114.02	122.70
1	A	338	VAL	N-CA-C	-5.42	96.37	111.00
1	B	296	ASP	CA-C-N	5.41	129.11	117.20
1	A	79	ARG	O-C-N	-5.40	114.06	122.70
1	B	131	CYS	O-C-N	-5.37	114.11	122.70
1	B	345	LEU	O-C-N	-5.32	114.18	122.70
1	A	179	GLN	N-CA-C	-5.22	96.91	111.00
1	B	427	VAL	CG1-CB-CG2	-5.21	102.57	110.90
1	A	235	VAL	CG1-CB-CG2	-5.20	102.57	110.90
1	A	345	LEU	O-C-N	-5.18	114.41	122.70
1	B	179	GLN	N-CA-C	-5.17	97.05	111.00
1	B	213	HIS	CA-C-N	-5.13	105.91	117.20
1	A	203	ASP	CA-C-N	-5.09	106.00	117.20
1	B	434	THR	N-CA-CB	5.07	119.94	110.30
1	B	235	VAL	CG1-CB-CG2	-5.05	102.81	110.90
1	A	241	SER	N-CA-C	-5.04	97.40	111.00
1	B	376	VAL	CG1-CB-CG2	5.01	118.92	110.90
1	B	295	GLN	N-CA-C	-5.01	97.48	111.00

There are no chirality outliers.

There are no planarity outliers.

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	3289	0	3120	93	0
1	B	3289	0	3121	110	12
2	A	43	0	30	7	0
2	B	43	0	30	24	0
3	A	5	0	0	1	0
3	B	5	0	0	2	0
4	A	2	0	0	0	0
4	B	2	0	0	0	0
5	A	4	0	0	0	0
5	B	4	0	0	0	0
All	All	6686	0	6301	185	12

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 14.

All (185) 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:434:THR:HG23	1:A:435:HIS:CB	1.24	1.65
1:A:434:THR:CG2	1:A:435:HIS:HB3	1.54	1.37
1:A:434:THR:CG2	1:A:435:HIS:CB	2.04	1.30
1:A:411:HIS:CE1	1:B:100:LYS:NZ	2.04	1.25
1:A:411:HIS:HE1	1:B:100:LYS:NZ	1.39	1.15
1:A:411:HIS:CE1	1:B:100:LYS:HZ2	1.60	1.12
1:B:99:GLU:HB3	1:B:100:LYS:HE2	1.34	1.10
1:A:99:GLU:HB3	1:A:100:LYS:HE2	1.34	1.09
1:A:413:ASP:OD2	1:B:99:GLU:OE2	1.69	1.09
1:B:434:THR:HG23	1:B:434:THR:O	1.50	1.04
1:A:411:HIS:NE2	1:B:99:GLU:OE1	1.91	1.03
1:A:434:THR:HG23	1:A:435:HIS:HB2	1.05	1.03
1:A:380:LEU:HA	3:A:501:PO4:O2	1.58	1.01

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:434:THR:HG23	1:A:435:HIS:HB3	1.00	0.99
1:B:380:LEU:HA	3:B:501:PO4:O2	1.67	0.94
1:A:434:THR:HG22	1:A:435:HIS:HB3	1.52	0.90
1:A:434:THR:CG2	1:A:435:HIS:HB2	1.87	0.90
1:B:213:HIS:HD2	2:B:500:HEM:C1D	1.90	0.89
1:A:385:PRO:HB3	1:A:390:SER:HB3	1.55	0.88
1:A:411:HIS:CE1	1:B:100:LYS:HZ3	1.90	0.88
1:B:204:TYR:HB3	2:B:500:HEM:C3A	2.10	0.87
1:B:214:ARG:HD2	2:B:500:HEM:C2C	2.10	0.86
1:B:213:HIS:CD2	2:B:500:HEM:C1D	2.65	0.85
1:B:434:THR:CG2	1:B:434:THR:O	2.17	0.83
1:B:385:PRO:HB3	1:B:390:SER:HB3	1.59	0.82
1:A:411:HIS:HE1	1:B:100:LYS:HZ2	0.89	0.81
1:B:380:LEU:HD23	1:B:386:LEU:HD22	1.62	0.80
1:B:213:HIS:HD2	2:B:500:HEM:ND	1.80	0.78
1:A:411:HIS:HE1	1:B:100:LYS:CE	1.96	0.78
1:B:213:HIS:HD2	2:B:500:HEM:C4D	2.02	0.78
1:A:411:HIS:CE1	1:B:100:LYS:CD	2.68	0.77
1:A:380:LEU:HD23	1:A:386:LEU:HD22	1.65	0.77
1:B:100:LYS:O	1:B:102:GLU:N	2.18	0.77
1:A:100:LYS:O	1:A:102:GLU:N	2.18	0.75
1:B:41:ASN:HD22	1:B:262:ARG:HH21	1.33	0.74
1:A:411:HIS:CE1	1:B:100:LYS:CE	2.70	0.74
1:A:413:ASP:CG	1:B:99:GLU:OE2	2.26	0.74
1:B:56:HIS:ND1	2:B:500:HEM:CBB	2.52	0.73
1:A:411:HIS:HE1	1:B:100:LYS:CD	2.03	0.72
1:B:94:TRP:HB3	1:B:96:TYR:CE2	2.25	0.71
1:B:214:ARG:HD2	2:B:500:HEM:C1C	2.26	0.71
1:B:213:HIS:CD2	2:B:500:HEM:C4D	2.79	0.71
1:A:411:HIS:CE1	1:B:99:GLU:OE1	2.45	0.69
1:A:94:TRP:HB3	1:A:96:TYR:CE2	2.27	0.68
1:A:82:HIS:CE1	1:A:262:ARG:HG2	2.28	0.68
1:A:413:ASP:HB2	1:B:99:GLU:OE2	1.94	0.68
1:B:183:PHE:HE2	2:B:500:HEM:CGA	2.07	0.68
1:A:428:SER:HB2	1:A:433:CYS:HB2	1.75	0.66
1:B:41:ASN:HD22	1:B:262:ARG:NH2	1.93	0.66
1:B:204:TYR:CD1	2:B:500:HEM:C1A	2.85	0.66
1:A:410:ARG:HG3	1:A:414:LYS:HD2	1.79	0.65
1:B:410:ARG:HG3	1:B:414:LYS:HD2	1.80	0.64
1:A:226:GLU:HG2	1:A:267:SER:H	1.63	0.63
1:A:413:ASP:CB	1:B:99:GLU:OE2	2.46	0.63

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:500:HEM:HMB2	2:B:500:HEM:HBB2	1.81	0.62
1:B:226:GLU:HG2	1:B:267:SER:H	1.64	0.62
1:A:413:ASP:HB2	1:B:99:GLU:CD	2.20	0.61
1:A:411:HIS:CE1	1:B:100:LYS:HD3	2.35	0.61
1:A:56:HIS:CE1	2:A:500:HEM:CBB	2.84	0.60
1:B:204:TYR:HB3	2:B:500:HEM:C2A	2.37	0.59
1:B:49:ASP:HA	1:B:73:PRO:HB3	1.85	0.58
1:A:49:ASP:HA	1:A:73:PRO:HB3	1.86	0.58
1:A:434:THR:CB	1:A:435:HIS:HB2	2.34	0.57
1:B:56:HIS:CE1	2:B:500:HEM:CBB	2.88	0.57
1:B:183:PHE:CE2	2:B:500:HEM:CGA	2.88	0.57
1:A:406:LEU:HD22	1:A:430:LEU:HD23	1.87	0.56
1:B:286:SER:HB3	1:B:291:LEU:HD23	1.86	0.56
1:A:238:ALA:HB3	1:A:250:PHE:HB2	1.87	0.56
1:B:434:THR:O	1:B:435:HIS:CB	2.39	0.56
1:A:91:ASP:HB3	1:A:118:PHE:CE2	2.41	0.56
1:B:72:GLY:HA3	1:B:89:LYS:NZ	2.21	0.56
1:B:91:ASP:HB3	1:B:118:PHE:CE2	2.40	0.55
1:B:238:ALA:HB3	1:B:250:PHE:HB2	1.88	0.55
1:A:428:SER:HB2	1:A:433:CYS:CB	2.35	0.55
1:A:123:ALA:HA	1:A:137:LEU:O	2.07	0.55
1:A:56:HIS:CE1	2:A:500:HEM:HBB1	2.42	0.55
1:A:286:SER:HB3	1:A:291:LEU:HD23	1.88	0.54
1:A:327:VAL:HG23	1:A:328:ILE:HD12	1.88	0.54
1:A:269:PRO:HD2	1:A:272:HIS:CD2	2.43	0.54
2:A:500:HEM:HBB2	2:A:500:HEM:HMB2	1.89	0.54
1:A:82:HIS:NE2	1:A:262:ARG:HG2	2.21	0.54
1:B:410:ARG:CG	1:B:414:LYS:HD2	2.38	0.54
1:B:183:PHE:HE2	2:B:500:HEM:O2A	1.91	0.54
2:A:500:HEM:HMC2	2:A:500:HEM:HBC2	1.89	0.54
1:A:56:HIS:HE1	2:A:500:HEM:HBB1	1.74	0.53
1:A:410:ARG:CG	1:A:414:LYS:HD2	2.37	0.53
1:B:269:PRO:HD2	1:B:272:HIS:CD2	2.45	0.52
1:B:327:VAL:HG23	1:B:328:ILE:HD12	1.91	0.52
1:A:392:SER:HB3	1:A:410:ARG:HA	1.92	0.52
1:B:406:LEU:HD22	1:B:430:LEU:HD23	1.92	0.52
1:A:131:CYS:SG	1:A:188:PRO:HB2	2.51	0.51
1:A:413:ASP:CB	1:B:99:GLU:CD	2.79	0.51
1:B:123:ALA:HA	1:B:137:LEU:O	2.11	0.51
1:B:204:TYR:HD1	2:B:500:HEM:C1A	2.29	0.51
1:B:392:SER:HB3	1:B:410:ARG:HA	1.93	0.51

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:176:TYR:HE1	2:B:500:HEM:O2A	1.94	0.50
1:B:214:ARG:HD2	2:B:500:HEM:C3C	2.46	0.50
1:A:294:ILE:HG12	1:A:299:VAL:HG22	1.92	0.50
1:B:328:ILE:HG22	1:B:329:SER:N	2.27	0.50
1:B:54:LYS:HZ1	1:B:66:ARG:HH21	1.59	0.50
1:B:216:SER:O	1:B:217:HIS:ND1	2.45	0.49
1:A:48:LYS:HB3	1:A:53:TRP:CZ3	2.47	0.49
1:B:412:VAL:O	1:B:415:LEU:HB3	2.12	0.49
1:B:434:THR:O	1:B:435:HIS:HB2	2.12	0.49
1:A:410:ARG:HG3	1:A:414:LYS:CD	2.43	0.48
1:B:294:ILE:HG12	1:B:299:VAL:HG22	1.94	0.48
1:A:216:SER:O	1:A:217:HIS:ND1	2.46	0.48
1:B:390:SER:HB2	1:B:397:ASN:ND2	2.29	0.47
1:B:333:VAL:HG21	1:B:347:ILE:HD11	1.96	0.47
1:A:379:ALA:HB2	1:A:400:LEU:HD23	1.94	0.47
1:A:328:ILE:HG22	1:A:329:SER:N	2.29	0.47
1:A:48:LYS:HB3	1:A:53:TRP:HZ3	1.79	0.47
1:B:379:ALA:HB2	1:B:400:LEU:HD23	1.95	0.47
1:B:24:ILE:HG23	1:B:25:GLU:H	1.78	0.47
1:B:48:LYS:HB3	1:B:53:TRP:CZ3	2.49	0.47
1:A:26:GLN:HE21	1:A:26:GLN:H	1.62	0.47
1:A:411:HIS:CD2	1:B:99:GLU:OE1	2.63	0.47
1:B:94:TRP:NE1	1:B:106:PRO:HG3	2.31	0.46
1:A:412:VAL:O	1:A:415:LEU:HB3	2.15	0.46
1:A:411:HIS:NE2	1:B:100:LYS:HD3	2.30	0.46
1:A:333:VAL:HG21	1:A:347:ILE:HD11	1.97	0.46
1:B:243:ASN:ND2	1:B:243:ASN:H	2.14	0.46
1:B:26:GLN:H	1:B:26:GLN:HE21	1.62	0.46
1:B:410:ARG:HG3	1:B:414:LYS:CD	2.44	0.46
1:A:122:ALA:HB3	1:A:139:PHE:HB2	1.98	0.46
1:A:24:ILE:HG23	1:A:25:GLU:H	1.79	0.46
1:A:327:VAL:HG11	1:B:389:ASN:HB3	1.98	0.46
1:A:390:SER:HB2	1:A:397:ASN:ND2	2.30	0.46
1:B:214:ARG:HG3	2:B:500:HEM:CB3	2.45	0.46
1:A:62:LEU:O	1:A:65:GLU:HB3	2.16	0.46
1:A:54:LYS:HE3	1:A:66:ARG:HH21	1.81	0.46
1:A:414:LYS:HG3	1:B:100:LYS:HD2	1.98	0.46
1:A:26:GLN:NE2	1:A:26:GLN:H	2.13	0.46
1:B:56:HIS:CD2	1:B:56:HIS:C	2.89	0.46
1:B:50:GLU:O	1:B:62:LEU:HD12	2.16	0.45
1:A:243:ASN:H	1:A:243:ASN:ND2	2.14	0.45

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:100:LYS:HA	1:B:100:LYS:HD3	1.44	0.45
1:B:426:ARG:HB3	1:B:426:ARG:HH11	1.81	0.45
1:B:122:ALA:HB3	1:B:139:PHE:HB2	1.99	0.45
1:B:131:CYS:SG	1:B:188:PRO:HB2	2.57	0.45
1:B:48:LYS:HB3	1:B:53:TRP:HZ3	1.81	0.45
1:B:100:LYS:HE2	1:B:100:LYS:N	2.32	0.45
1:B:62:LEU:O	1:B:65:GLU:HB3	2.17	0.44
1:A:100:LYS:N	1:A:100:LYS:HE2	2.32	0.44
1:A:50:GLU:O	1:A:62:LEU:HD12	2.18	0.44
1:B:390:SER:HB2	1:B:397:ASN:HD21	1.82	0.44
1:A:399:TYR:HA	1:A:407:TYR:O	2.16	0.44
2:A:500:HEM:HBD2	2:A:500:HEM:HHA	1.99	0.44
1:B:214:ARG:CD	2:B:500:HEM:C3C	3.00	0.44
1:B:55:SER:OG	1:B:57:ARG:HB2	2.18	0.44
1:B:72:GLY:HA3	1:B:89:LYS:HZ2	1.82	0.44
1:A:72:GLY:HA3	1:A:89:LYS:NZ	2.33	0.43
1:B:223:HIS:CD2	1:B:234:LEU:HD12	2.52	0.43
1:B:204:TYR:CD1	2:B:500:HEM:C2A	3.06	0.43
1:A:56:HIS:ND1	2:A:500:HEM:CBB	2.81	0.43
1:A:214:ARG:HD3	1:A:214:ARG:N	2.33	0.43
1:B:258:LEU:HD23	1:B:265:TRP:HZ3	1.82	0.43
1:A:56:HIS:CD2	1:A:56:HIS:C	2.92	0.43
1:A:55:SER:OG	1:A:57:ARG:HB2	2.19	0.43
1:B:214:ARG:HD3	1:B:214:ARG:N	2.33	0.43
1:B:26:GLN:NE2	1:B:26:GLN:H	2.15	0.43
1:A:174:ARG:HH22	1:A:272:HIS:CD2	2.37	0.43
1:B:386:LEU:HG	1:B:430:LEU:HD22	2.01	0.43
1:B:399:TYR:HA	1:B:407:TYR:O	2.18	0.42
1:B:204:TYR:CD1	2:B:500:HEM:CHA	3.02	0.42
1:A:426:ARG:HB3	1:A:426:ARG:HH11	1.84	0.42
1:A:386:LEU:HG	1:A:430:LEU:HD22	2.01	0.42
1:A:54:LYS:CE	1:A:66:ARG:HH21	2.33	0.42
1:A:223:HIS:CD2	1:A:234:LEU:HD12	2.54	0.42
1:B:24:ILE:HG23	1:B:25:GLU:N	2.34	0.42
1:A:274:TRP:HA	1:A:275:PRO:HD2	1.86	0.42
1:B:386:LEU:HA	1:B:386:LEU:HD12	1.78	0.42
1:A:344:ARG:HG2	1:A:355:TRP:CE3	2.55	0.42
1:A:346:HIS:HA	1:A:354:TRP:O	2.20	0.42
1:A:99:GLU:HB3	1:A:100:LYS:CE	2.25	0.42
1:A:100:LYS:HD3	1:A:100:LYS:HA	1.44	0.41
1:B:54:LYS:NZ	1:B:66:ARG:HH21	2.17	0.41

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:82:HIS:ND1	1:B:82:HIS:O	2.53	0.41
1:B:256:TRP:CE2	1:B:267:SER:OG	2.72	0.41
1:A:258:LEU:HD23	1:A:265:TRP:HZ3	1.85	0.41
1:B:94:TRP:CE2	1:B:106:PRO:HG3	2.55	0.41
1:B:213:HIS:CD2	2:B:500:HEM:CHD	3.03	0.41
1:B:285:PHE:HA	3:B:501:PO4:O4	2.20	0.41
1:B:274:TRP:HA	1:B:275:PRO:HD2	1.85	0.40
1:A:100:LYS:C	1:A:102:GLU:N	2.75	0.40

All (12) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:118:PHE:CZ	1:B:312:ASN:CG[4_556]	1.31	0.89
1:B:51:PHE:CE1	1:B:156:GLU:O[4_556]	1.44	0.76
1:B:118:PHE:CE1	1:B:312:ASN:ND2[4_556]	1.73	0.47
1:B:118:PHE:CZ	1:B:312:ASN:OD1[4_556]	1.80	0.40
1:B:118:PHE:CE2	1:B:312:ASN:CB[4_556]	1.88	0.32
1:B:118:PHE:CZ	1:B:312:ASN:CB[4_556]	1.92	0.28
1:B:118:PHE:CZ	1:B:312:ASN:ND2[4_556]	1.93	0.27
1:B:25:GLU:OE2	1:B:154:LYS:O[4_556]	1.98	0.22
1:B:118:PHE:CE1	1:B:312:ASN:CG[4_556]	2.11	0.09
1:B:51:PHE:CD1	1:B:156:GLU:O[4_556]	2.14	0.06
1:B:60:ARG:NH2	1:B:156:GLU:OE1[4_556]	2.16	0.04
1:B:62:LEU:CD1	1:B:158:SER:CB[4_556]	2.18	0.02

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	404/460 (88%)	370 (92%)	30 (7%)	4 (1%)	15 45

Continued on next page...

Continued from previous page...

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	B	404/460 (88%)	369 (91%)	31 (8%)	4 (1%)	15	45
All	All	808/920 (88%)	739 (92%)	61 (8%)	8 (1%)	15	45

All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	101	ASN
1	A	296	ASP
1	B	101	ASN
1	B	296	ASP
1	A	55	SER
1	B	55	SER
1	B	82	HIS
1	A	82	HIS

5.3.2 Protein sidechains ⓘ

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	352/393 (90%)	317 (90%)	35 (10%)	8	24
1	B	352/393 (90%)	318 (90%)	34 (10%)	8	25
All	All	704/786 (90%)	635 (90%)	69 (10%)	8	24

All (69) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	26	GLN
1	A	43	THR
1	A	55	SER
1	A	56	HIS
1	A	66	ARG
1	A	80	HIS
1	A	92	LYS
1	A	98	SER

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	A	99	GLU
1	A	100	LYS
1	A	108	SER
1	A	177	CYS
1	A	206	LEU
1	A	214	ARG
1	A	217	HIS
1	A	218	ARG
1	A	227	SER
1	A	228	THR
1	A	240	VAL
1	A	249	VAL
1	A	256	TRP
1	A	272	HIS
1	A	279	SER
1	A	305	LYS
1	A	327	VAL
1	A	329	SER
1	A	344	ARG
1	A	356	LEU
1	A	371	TRP
1	A	376	VAL
1	A	394	SER
1	A	410	ARG
1	A	411	HIS
1	A	426	ARG
1	A	429	ARG
1	B	26	GLN
1	B	43	THR
1	B	55	SER
1	B	56	HIS
1	B	66	ARG
1	B	80	HIS
1	B	92	LYS
1	B	98	SER
1	B	99	GLU
1	B	100	LYS
1	B	108	SER
1	B	177	CYS
1	B	206	LEU
1	B	214	ARG
1	B	217	HIS

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	B	218	ARG
1	B	227	SER
1	B	228	THR
1	B	240	VAL
1	B	249	VAL
1	B	256	TRP
1	B	279	SER
1	B	305	LYS
1	B	327	VAL
1	B	329	SER
1	B	344	ARG
1	B	356	LEU
1	B	371	TRP
1	B	376	VAL
1	B	394	SER
1	B	410	ARG
1	B	411	HIS
1	B	426	ARG
1	B	429	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (12) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	26	GLN
1	A	41	ASN
1	A	56	HIS
1	A	142	ASN
1	A	243	ASN
1	A	411	HIS
1	B	26	GLN
1	B	41	ASN
1	B	142	ASN
1	B	243	ASN
1	B	272	HIS
1	B	295	GLN

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 carbohydrates in this entry.

5.6 Ligand geometry ⓘ

Of 16 ligands modelled in this entry, 12 are monoatomic - leaving 4 for Mogul analysis.

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$
2	HEM	B	500	1	27,50,50	1.51	4 (14%)	17,82,82	3.58	11 (64%)
2	HEM	A	500	1	27,50,50	1.62	4 (14%)	17,82,82	1.94	6 (35%)
3	PO4	A	501	-	4,4,4	1.77	2 (50%)	6,6,6	1.55	1 (16%)
3	PO4	B	501	-	4,4,4	1.77	2 (50%)	6,6,6	1.55	1 (16%)

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	B	500	1	-	5/6/54/54	-
2	HEM	A	500	1	-	0/6/54/54	-

All (12) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	500	HEM	C3B-CAB	-5.70	1.36	1.47
2	B	500	HEM	CBC-CAC	4.25	1.57	1.29
2	A	500	HEM	C3C-CAC	-3.41	1.40	1.47

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	500	HEM	CBC-CAC	2.82	1.48	1.29
2	B	500	HEM	CBB-CAB	2.65	1.46	1.29
3	A	501	PO4	P-O2	-2.42	1.47	1.54
3	B	501	PO4	P-O2	-2.40	1.47	1.54
3	A	501	PO4	P-O4	-2.16	1.48	1.54
3	B	501	PO4	P-O4	-2.15	1.48	1.54
2	B	500	HEM	C1B-C2B	-2.11	1.37	1.42
2	A	500	HEM	C1A-NA	-2.04	1.32	1.36
2	B	500	HEM	C1A-NA	-2.01	1.32	1.36

All (19) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	500	HEM	CBD-CAD-C3D	7.13	125.61	112.48
2	B	500	HEM	CAD-CBD-CGD	6.41	123.42	112.67
2	B	500	HEM	CBA-CAA-C2A	5.73	123.05	112.49
2	A	500	HEM	C4A-C3A-C2A	-4.31	104.00	107.00
2	B	500	HEM	C4A-C3A-C2A	-4.25	104.04	107.00
2	B	500	HEM	C3B-C4B-NB	3.90	114.25	109.21
2	B	500	HEM	C1D-C2D-C3D	3.82	109.66	107.00
2	B	500	HEM	CMD-C2D-C3D	-3.42	118.49	124.94
2	B	500	HEM	CMA-C3A-C2A	3.22	131.01	124.94
3	B	501	PO4	O4-P-O1	3.09	122.20	110.89
3	A	501	PO4	O4-P-O1	3.08	122.17	110.89
2	A	500	HEM	CBA-CAA-C2A	2.89	117.82	112.49
2	A	500	HEM	C4C-C3C-C2C	-2.86	104.90	106.90
2	B	500	HEM	C3C-C4C-NC	-2.55	106.12	110.94
2	B	500	HEM	CMA-C3A-C4A	-2.29	124.95	128.46
2	A	500	HEM	CMB-C2B-C3B	2.26	128.91	124.68
2	A	500	HEM	CMA-C3A-C2A	2.13	128.96	124.94
2	B	500	HEM	CMB-C2B-C3B	2.05	128.51	124.68
2	A	500	HEM	CMC-C2C-C3C	2.00	128.42	124.68

There are no chirality outliers.

All (5) torsion outliers are listed below:

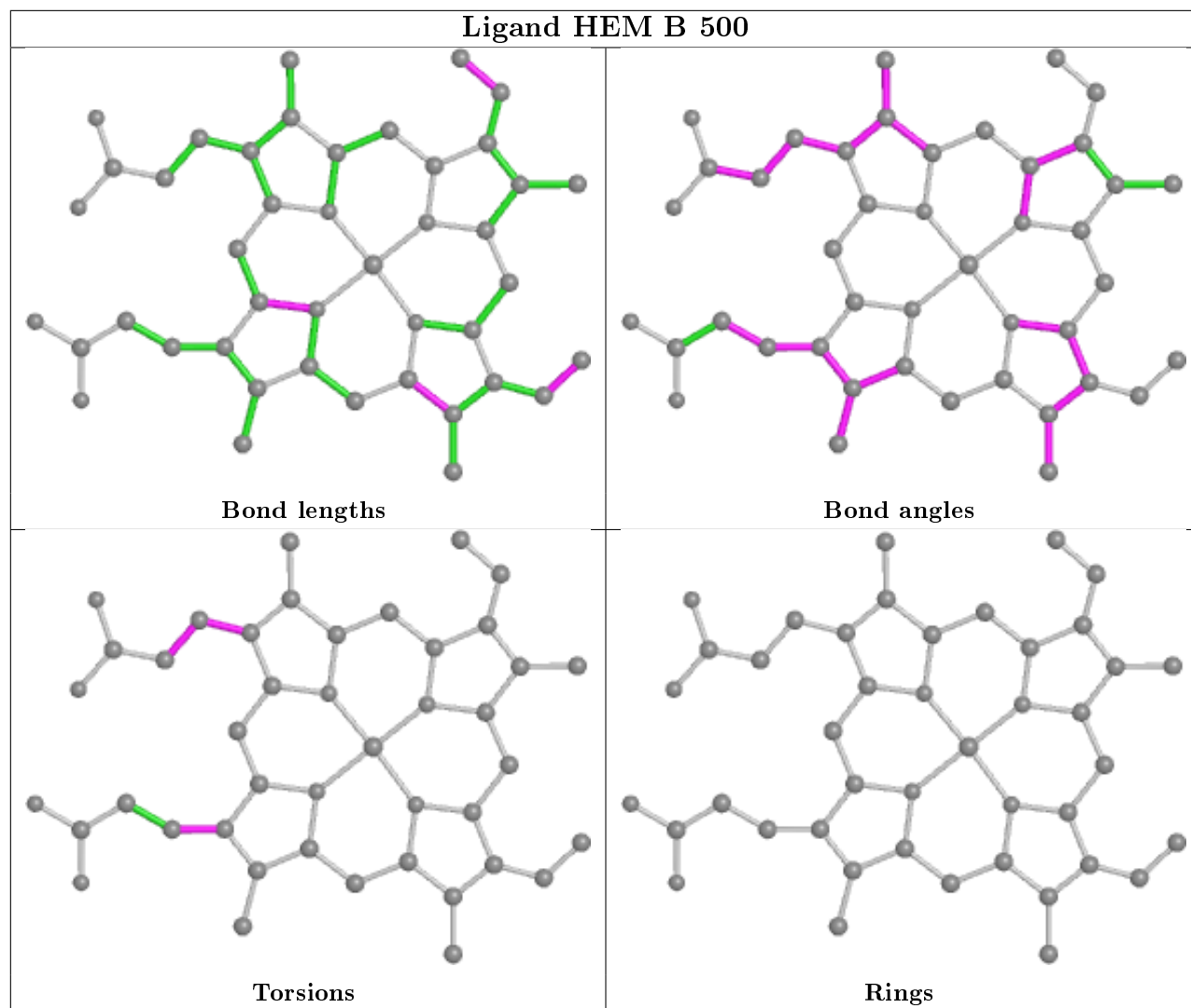
Mol	Chain	Res	Type	Atoms
2	B	500	HEM	C1A-C2A-CAA-CBA
2	B	500	HEM	C3A-C2A-CAA-CBA
2	B	500	HEM	C2D-C3D-CAD-CBD
2	B	500	HEM	C4D-C3D-CAD-CBD
2	B	500	HEM	C3D-CAD-CBD-CGD

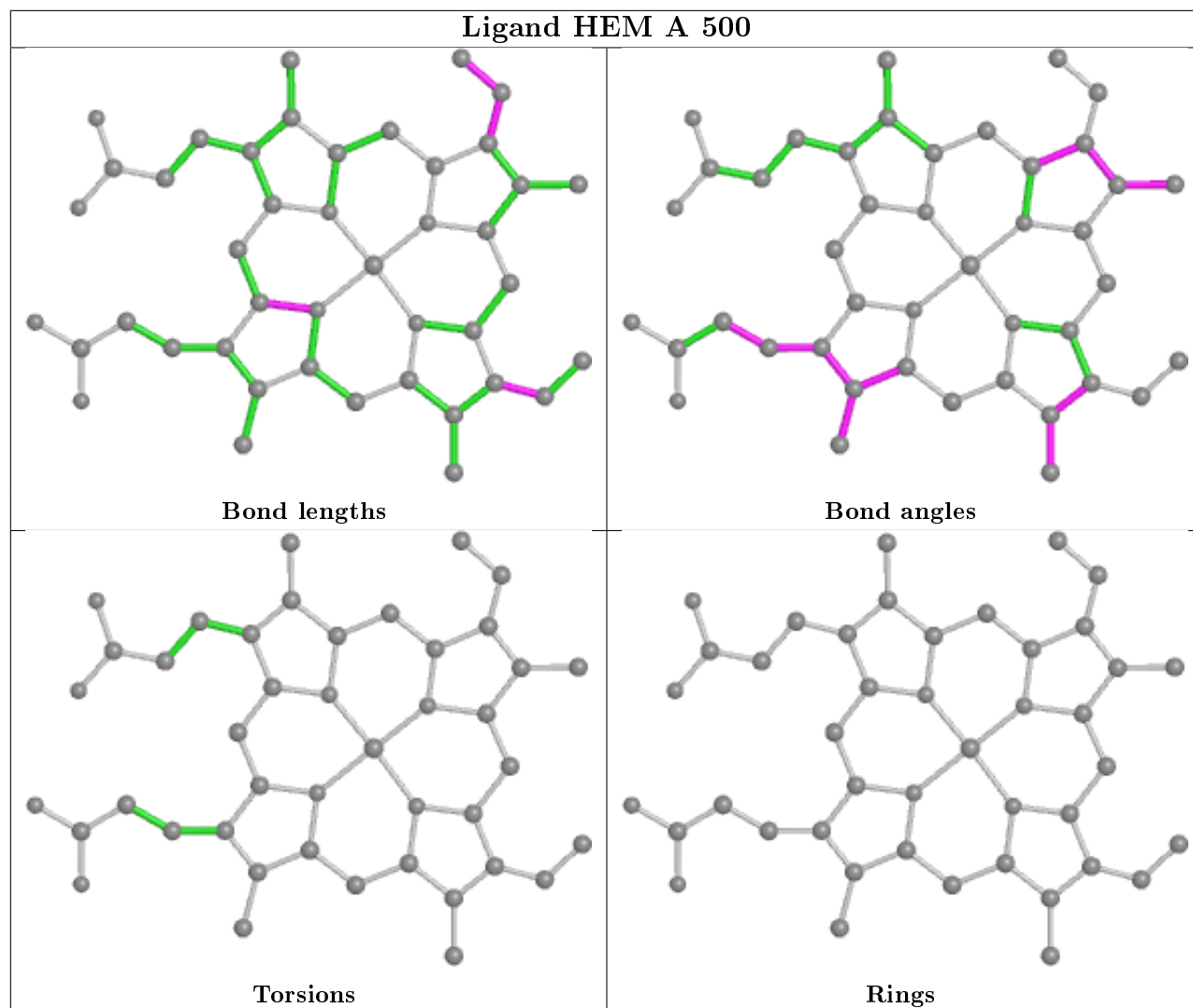
There are no ring outliers.

4 monomers are involved in 34 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	500	HEM	24	0
2	A	500	HEM	7	0
3	A	501	PO4	1	0
3	B	501	PO4	2	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](#)

The following chains have linkage breaks:

Mol	Chain	Number of breaks
1	B	1
1	A	1

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	B	434:THR	C	435:HIS	N	1.20
1	A	434:THR	C	435:HIS	N	1.13

6 Fit of model and data

6.1 Protein, DNA and RNA chains

Unable to reproduce the depositors R factor - this section is therefore empty.

6.2 Non-standard residues in protein, DNA, RNA chains

Unable to reproduce the depositors R factor - this section is therefore empty.

6.3 Carbohydrates

Unable to reproduce the depositors R factor - this section is therefore empty.

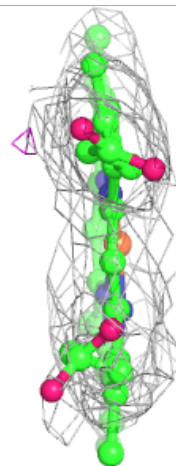
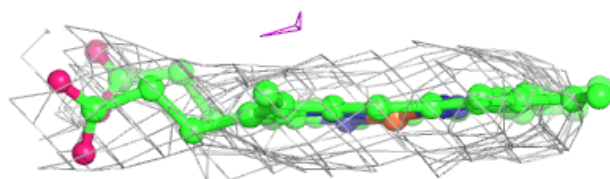
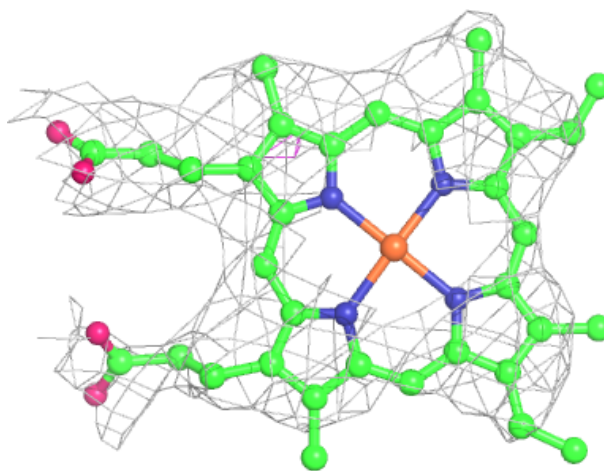
6.4 Ligands

Unable to reproduce the depositors R factor - this section is therefore empty.

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.

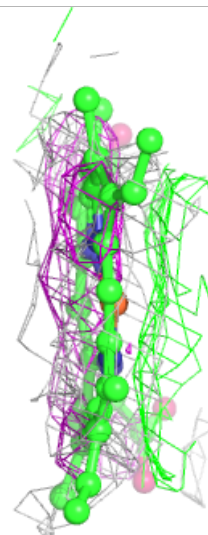
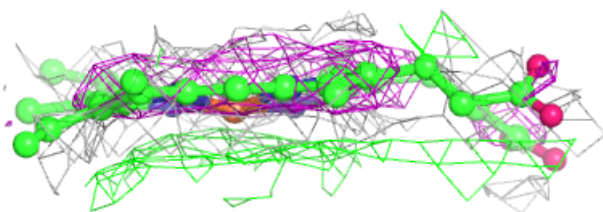
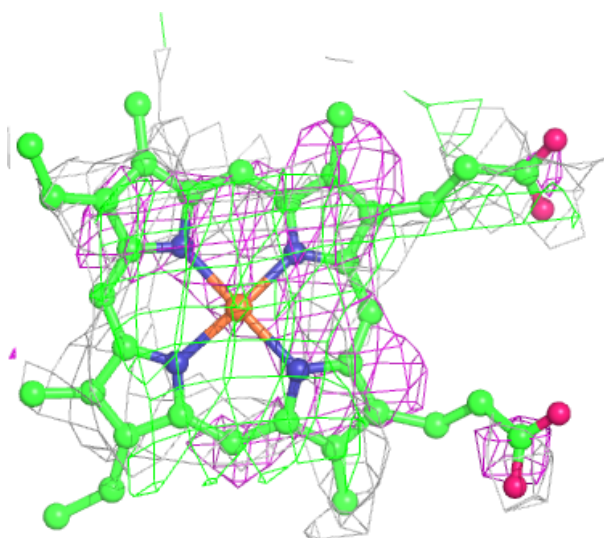
Electron density around HEM A 500:

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



Electron density around HEM B 500:

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



6.5 Other polymers [i](#)

Unable to reproduce the depositors R factor - this section is therefore empty.