Full wwPDB X-ray Structure Validation Report ⓘ

Oct 16, 2023 – 07:06 PM EDT

PDB ID	:	1S0H
Title	:	Structure determination of haemoglobin from Donkey(equus asinus) at 3.0 Angstrom resolution
Authors	:	Balasundaresan, D.; Ponnuswamy, M.N.; Saraboji, K.
Deposited on	:	2003-12-31
Resolution	:	3.00 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:

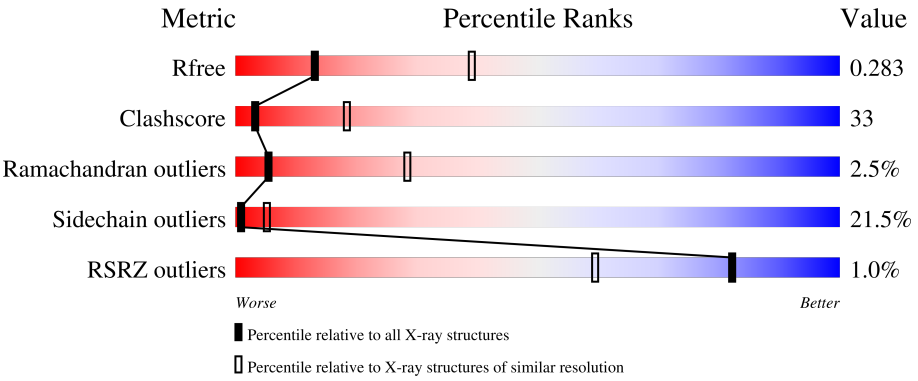
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.36
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 3.00 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}	130704	2092 (3.00-3.00)
Clashscore	141614	2416 (3.00-3.00)
Ramachandran outliers	138981	2333 (3.00-3.00)
Sidechain outliers	138945	2336 (3.00-3.00)
RSRZ outliers	127900	1990 (3.00-3.00)

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 <=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	141	41% 43% 14%
2	B	146	34% 47% 16%

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	HEM	B	147	-	-	X	-

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 2286 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.

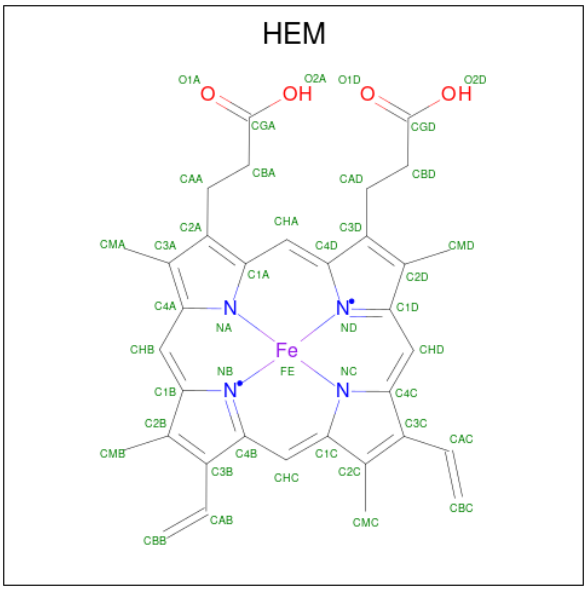
- Molecule 1 is a protein called Hemoglobin alpha chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	141	Total	C	N	O	S	0	0	0
			1066	682	186	196	2			

- Molecule 2 is a protein called Hemoglobin beta chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	146	Total	C	N	O	S	0	0	0
			1134	728	200	204	2			

- Molecule 3 is PROTOPORPHYRIN IX CONTAINING FE (three-letter code: HEM) (formula: $C_{34}H_{32}FeN_4O_4$).

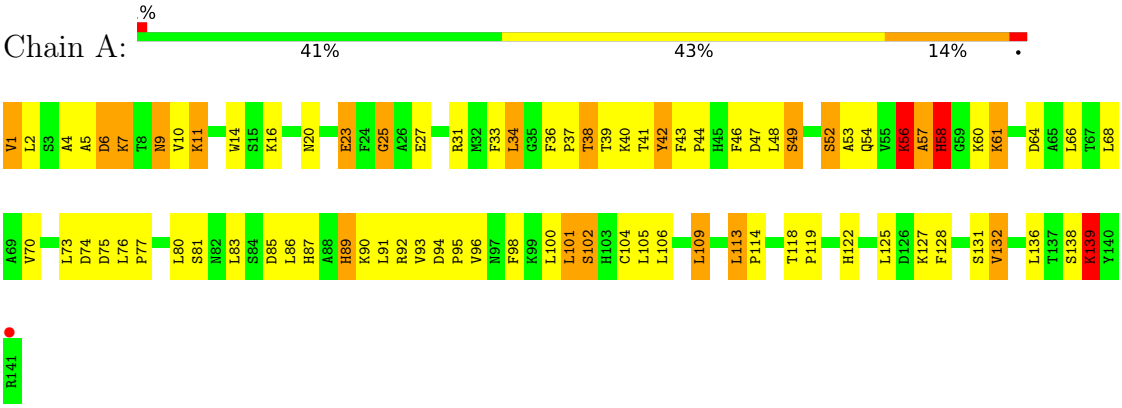


Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	Fe	N	O	0	0
			43	34	1	4	4		
3	B	1	Total	C	Fe	N	O	0	0
			43	34	1	4	4		

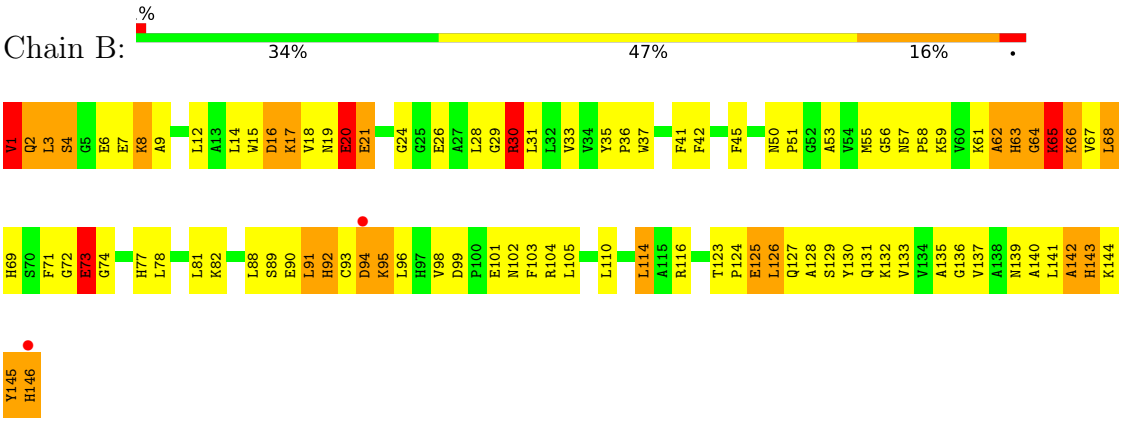
3 Residue-property plots (i)

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: Hemoglobin alpha chain



• Molecule 2: Hemoglobin beta chain



4 Data and refinement statistics (i)

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α, β, γ	107.66Å 63.08Å 54.04Å 90.00° 111.75° 90.00°	Depositor
Resolution (Å)	21.18 – 3.00 21.16 – 3.00	Depositor EDS
% Data completeness (in resolution range)	79.0 (21.18-3.00) 79.0 (21.16-3.00)	Depositor EDS
R_{merge}	0.16	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	1.04 (at 2.98Å)	Xtriage
Refinement program	REFMAC 5.1.24	Depositor
R, R_{free}	0.181 , 0.285 0.193 , 0.283	Depositor DCC
R_{free} test set	254 reflections (4.72%)	wwPDB-VP
Wilson B-factor (Å^2)	53.0	Xtriage
Anisotropy	0.163	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.33 , 84.0	EDS
L-test for twinning[2]	$< \|L\| > = 0.46$, $< L^2 > = 0.28$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.92	EDS
Total number of atoms	2286	wwPDB-VP
Average B, all atoms (Å^2)	34.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.43% of the height of the origin peak. No significant pseudotranslation is detected.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality (i)

5.1 Standard geometry (i)

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

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	1.86	16/1092 (1.5%)	1.62	15/1481 (1.0%)
2	B	1.98	27/1162 (2.3%)	1.95	24/1572 (1.5%)
All	All	1.92	43/2254 (1.9%)	1.80	39/3053 (1.3%)

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	2
2	B	0	3
All	All	0	5

All (43) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	58	HIS	C-N	-12.64	1.10	1.33
1	A	57	ALA	C-N	-10.74	1.09	1.34
2	B	91	LEU	C-N	9.28	1.55	1.34
2	B	59	LYS	CD-CE	7.74	1.70	1.51
1	A	127	LYS	CD-CE	7.63	1.70	1.51
2	B	92	HIS	C-N	7.35	1.50	1.34
2	B	30	ARG	CG-CD	-7.08	1.34	1.51
1	A	10	VAL	CB-CG2	7.05	1.67	1.52
2	B	18	VAL	CB-CG1	6.92	1.67	1.52
1	A	33	PHE	C-N	6.88	1.49	1.34
2	B	21	GLU	CD-OE2	6.87	1.33	1.25
1	A	90	LYS	CB-CG	6.61	1.70	1.52
2	B	101	GLU	CD-OE2	6.59	1.32	1.25
2	B	130	TYR	CD2-CE2	6.47	1.49	1.39

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	11	LYS	CD-CE	6.46	1.67	1.51
2	B	130	TYR	CB-CG	-6.42	1.42	1.51
2	B	18	VAL	CB-CG2	6.36	1.66	1.52
1	A	90	LYS	CD-CE	6.28	1.67	1.51
2	B	101	GLU	CG-CD	6.28	1.61	1.51
1	A	132	VAL	CB-CG1	6.27	1.66	1.52
2	B	35	TYR	CE1-CZ	-6.24	1.30	1.38
2	B	73	GLU	CG-CD	6.17	1.61	1.51
2	B	20	GLU	CD-OE1	6.00	1.32	1.25
2	B	74	GLY	C-O	-5.99	1.14	1.23
1	A	14	TRP	CE3-CZ3	-5.81	1.28	1.38
2	B	143	HIS	C-O	5.78	1.34	1.23
2	B	114	LEU	C-O	-5.76	1.12	1.23
1	A	14	TRP	CB-CG	-5.58	1.40	1.50
2	B	66	LYS	CE-NZ	5.55	1.62	1.49
2	B	35	TYR	CG-CD1	-5.49	1.32	1.39
1	A	118	THR	C-O	-5.39	1.13	1.23
2	B	63	HIS	C-N	-5.39	1.23	1.33
2	B	66	LYS	CB-CG	5.38	1.67	1.52
2	B	37	TRP	CE3-CZ3	-5.34	1.29	1.38
1	A	42	TYR	CE2-CZ	5.30	1.45	1.38
2	B	142	ALA	CA-C	5.29	1.66	1.52
2	B	17	LYS	CE-NZ	5.22	1.62	1.49
1	A	128	PHE	CB-CG	-5.21	1.42	1.51
2	B	2	GLN	C-O	5.15	1.33	1.23
2	B	62	ALA	C-N	-5.06	1.22	1.34
2	B	30	ARG	CZ-NH2	-5.04	1.26	1.33
1	A	14	TRP	CG-CD1	-5.03	1.29	1.36
1	A	56	LYS	CD-CE	5.03	1.63	1.51

All (39) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	B	63	HIS	O-C-N	-21.67	86.36	123.20
2	B	63	HIS	CA-C-N	18.17	152.54	116.20
2	B	62	ALA	O-C-N	14.90	146.53	122.70
2	B	91	LEU	O-C-N	14.29	145.57	122.70
1	A	57	ALA	O-C-N	13.99	145.09	122.70
2	B	63	HIS	C-N-CA	12.73	149.04	122.30
2	B	92	HIS	O-C-N	-12.20	103.19	122.70
2	B	62	ALA	CA-C-N	-11.45	92.00	117.20
2	B	91	LEU	CA-C-N	-10.79	93.45	117.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	58	HIS	O-C-N	-9.57	106.92	123.20
1	A	57	ALA	CA-C-N	-9.36	96.61	117.20
2	B	92	HIS	C-N-CA	9.20	144.71	121.70
2	B	65	LYS	C-N-CA	-8.66	100.06	121.70
2	B	92	HIS	CA-C-N	8.23	135.31	117.20
2	B	62	ALA	C-N-CA	-7.50	102.95	121.70
1	A	52	SER	CA-C-N	-7.46	100.80	117.20
1	A	58	HIS	CA-C-N	6.68	129.55	116.20
2	B	99	ASP	CB-CG-OD2	5.98	123.68	118.30
1	A	52	SER	O-C-N	5.94	132.21	122.70
1	A	57	ALA	C-N-CA	-5.90	106.95	121.70
2	B	65	LYS	O-C-N	5.71	131.83	122.70
1	A	139	LYS	CD-CE-NZ	5.67	124.73	111.70
2	B	2	GLN	CA-C-N	-5.66	104.75	117.20
1	A	6	ASP	CB-CG-OD2	5.62	123.36	118.30
2	B	30	ARG	CG-CD-NE	5.52	123.39	111.80
2	B	16	ASP	CB-CG-OD2	5.52	123.26	118.30
1	A	80	LEU	CB-CG-CD1	5.43	120.23	111.00
1	A	106	LEU	CB-CG-CD1	-5.40	101.82	111.00
2	B	125	GLU	N-CA-C	-5.33	96.62	111.00
2	B	146	HIS	CA-C-O	5.33	131.29	120.10
2	B	59	LYS	CD-CE-NZ	5.30	123.88	111.70
1	A	68	LEU	CB-CG-CD1	5.29	119.99	111.00
1	A	74	ASP	CB-CG-OD2	5.23	123.01	118.30
1	A	34	LEU	O-C-N	-5.23	114.31	123.20
2	B	2	GLN	CB-CA-C	-5.23	99.94	110.40
1	A	25	GLY	N-CA-C	5.21	126.12	113.10
2	B	2	GLN	O-C-N	5.13	130.92	122.70
2	B	36	PRO	N-CD-CG	-5.11	95.53	103.20
2	B	30	ARG	CD-NE-CZ	5.10	130.73	123.60

There are no chirality outliers.

All (5) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	49	SER	Peptide
1	A	58	HIS	Mainchain
2	B	1	VAL	Peptide
2	B	3	LEU	Peptide
2	B	89	SER	Peptide

5.2 Too-close contacts (i)

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	1066	0	1070	53	0
2	B	1134	0	1123	99	0
3	A	43	0	30	3	0
3	B	43	0	30	27	0
All	All	2286	0	2253	150	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 33.

All (150) 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)
2:B:88:LEU:HD21	3:B:147:HEM:HBA1	1.26	1.13
2:B:141:LEU:HD11	3:B:147:HEM:HMB3	1.36	1.07
2:B:88:LEU:CD2	3:B:147:HEM:HBA1	1.89	1.01
2:B:88:LEU:HD11	3:B:147:HEM:CMA	1.94	0.97
1:A:4:ALA:HA	1:A:7:LYS:HG3	1.50	0.93
2:B:88:LEU:HD11	3:B:147:HEM:HMA2	1.49	0.91
2:B:142:ALA:O	2:B:145:TYR:HB2	1.72	0.89
2:B:141:LEU:HD11	3:B:147:HEM:CMB	2.02	0.89
1:A:37:PRO:HB3	1:A:40:LYS:HE3	1.60	0.81
2:B:88:LEU:CD1	3:B:147:HEM:HMA1	2.10	0.81
2:B:88:LEU:CD1	3:B:147:HEM:CMA	2.60	0.80
1:A:43:PHE:N	1:A:44:PRO:HD3	1.97	0.79
1:A:42:TYR:C	1:A:44:PRO:HD3	2.07	0.75
1:A:70:VAL:O	1:A:73:LEU:HG	1.89	0.73
2:B:141:LEU:CD1	3:B:147:HEM:HMB3	2.17	0.72
2:B:20:GLU:OE2	2:B:65:LYS:NZ	2.23	0.72
1:A:37:PRO:HB3	1:A:40:LYS:CE	2.21	0.69
2:B:88:LEU:HD11	3:B:147:HEM:HMA1	1.72	0.68
2:B:65:LYS:O	2:B:66:LYS:C	2.29	0.67
1:A:1:VAL:HG21	1:A:131:SER:OG	1.96	0.65
2:B:42:PHE:CE2	3:B:147:HEM:HMD1	2.32	0.65
2:B:103:PHE:HB3	3:B:147:HEM:HAB	1.79	0.65
2:B:126:LEU:O	2:B:126:LEU:HD23	1.97	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:47:ASP:O	1:A:52:SER:CB	2.46	0.64
2:B:141:LEU:HD11	3:B:147:HEM:C2B	2.33	0.63
2:B:110:LEU:O	2:B:110:LEU:HD23	1.98	0.63
2:B:126:LEU:O	2:B:126:LEU:CD2	2.46	0.63
2:B:73:GLU:O	2:B:77:HIS:HB2	2.00	0.62
2:B:88:LEU:HD13	3:B:147:HEM:HMA1	1.81	0.62
2:B:131:GLN:NE2	2:B:131:GLN:HA	2.16	0.61
1:A:75:ASP:CG	1:A:75:ASP:O	2.37	0.60
2:B:131:GLN:HE21	2:B:131:GLN:CA	2.12	0.60
2:B:131:GLN:HA	2:B:131:GLN:HE21	1.67	0.60
1:A:76:LEU:N	1:A:77:PRO:CD	2.65	0.60
2:B:66:LYS:O	2:B:69:HIS:HB3	2.03	0.59
1:A:43:PHE:N	1:A:44:PRO:CD	2.66	0.59
2:B:92:HIS:CE1	3:B:147:HEM:NA	2.70	0.58
2:B:41:PHE:CD1	3:B:147:HEM:HBC1	2.39	0.58
1:A:46:PHE:HA	1:A:54:GLN:OE1	2.05	0.57
1:A:37:PRO:CB	1:A:40:LYS:HE3	2.33	0.56
1:A:113:LEU:N	1:A:114:PRO:HD3	2.21	0.56
2:B:91:LEU:HD22	3:B:147:HEM:HBA2	1.88	0.56
2:B:9:ALA:HA	2:B:12:LEU:HD12	1.87	0.56
1:A:101:LEU:O	1:A:102:SER:C	2.43	0.55
2:B:24:GLY:HA3	2:B:64:GLY:O	2.05	0.55
2:B:63:HIS:CE1	3:B:147:HEM:HAD2	2.41	0.55
2:B:14:LEU:O	2:B:17:LYS:N	2.40	0.55
1:A:47:ASP:O	1:A:52:SER:HB2	2.07	0.55
2:B:3:LEU:HD12	2:B:8:LYS:CE	2.37	0.55
2:B:8:LYS:O	2:B:12:LEU:HD12	2.06	0.55
2:B:129:SER:O	2:B:132:LYS:HB2	2.07	0.55
2:B:3:LEU:HD12	2:B:8:LYS:HE2	1.89	0.54
1:A:2:LEU:HD22	1:A:6:ASP:HB3	1.90	0.53
2:B:110:LEU:HD23	2:B:110:LEU:C	2.28	0.53
2:B:24:GLY:HA2	2:B:68:LEU:HD23	1.90	0.53
1:A:100:LEU:O	1:A:101:LEU:C	2.42	0.53
2:B:131:GLN:NE2	2:B:131:GLN:CA	2.72	0.53
2:B:141:LEU:CD1	3:B:147:HEM:CMB	2.79	0.53
2:B:1:VAL:N	2:B:78:LEU:O	2.42	0.52
1:A:54:GLN:O	1:A:57:ALA:HB3	2.09	0.52
2:B:110:LEU:HD22	2:B:114:LEU:HD12	1.90	0.52
2:B:110:LEU:C	2:B:110:LEU:CD2	2.77	0.52
2:B:50:ASN:HD21	2:B:53:ALA:HB2	1.74	0.51
1:A:5:ALA:O	1:A:9:ASN:HB2	2.10	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:61:LYS:O	1:A:64:ASP:HB2	2.12	0.50
1:A:37:PRO:C	1:A:39:THR:N	2.63	0.50
2:B:4:SER:HB3	2:B:7:GLU:CD	2.32	0.50
2:B:50:ASN:HB2	2:B:51:PRO:HD2	1.93	0.50
2:B:90:GLU:O	2:B:91:LEU:C	2.50	0.50
2:B:95:LYS:HZ1	2:B:96:LEU:N	2.10	0.50
2:B:135:ALA:O	2:B:136:GLY:C	2.50	0.49
1:A:119:PRO:O	1:A:122:HIS:HB3	2.12	0.49
1:A:87:HIS:CE1	3:A:142:HEM:NA	2.78	0.49
1:A:66:LEU:O	1:A:70:VAL:HG23	2.13	0.48
2:B:14:LEU:C	2:B:16:ASP:N	2.64	0.48
2:B:129:SER:O	2:B:133:VAL:HG23	2.13	0.48
2:B:143:HIS:C	2:B:145:TYR:H	2.15	0.48
2:B:103:PHE:CB	3:B:147:HEM:HAB	2.44	0.48
1:A:89:HIS:HB2	1:A:139:LYS:CE	2.44	0.47
1:A:122:HIS:CD2	1:A:122:HIS:C	2.88	0.47
1:A:93:VAL:HG11	3:A:142:HEM:CAC	2.44	0.47
2:B:63:HIS:CE1	3:B:147:HEM:C3D	3.03	0.47
2:B:95:LYS:HZ1	2:B:95:LYS:C	2.18	0.47
1:A:131:SER:O	1:A:132:VAL:C	2.49	0.47
2:B:57:ASN:HA	2:B:58:PRO:HD3	1.64	0.47
1:A:37:PRO:C	1:A:39:THR:H	2.18	0.47
2:B:30:ARG:O	2:B:31:LEU:C	2.51	0.47
2:B:19:ASN:OD1	2:B:19:ASN:C	2.53	0.47
2:B:93:CYS:HB2	2:B:145:TYR:CE2	2.50	0.46
1:A:76:LEU:N	1:A:77:PRO:HD3	2.30	0.46
2:B:110:LEU:HD22	2:B:114:LEU:CD1	2.46	0.46
2:B:15:TRP:CZ2	2:B:72:GLY:HA2	2.50	0.46
1:A:20:ASN:O	1:A:23:GLU:N	2.49	0.45
2:B:95:LYS:C	2:B:95:LYS:NZ	2.69	0.45
2:B:8:LYS:O	2:B:12:LEU:CD1	2.64	0.45
2:B:98:VAL:O	2:B:145:TYR:OH	2.34	0.45
2:B:102:ASN:O	2:B:105:LEU:HB2	2.16	0.45
2:B:26:GLU:O	2:B:30:ARG:HD3	2.16	0.45
2:B:92:HIS:CE1	3:B:147:HEM:C4A	3.05	0.45
2:B:98:VAL:HG11	2:B:103:PHE:CE1	2.52	0.45
2:B:63:HIS:HE1	3:B:147:HEM:HAD2	1.80	0.45
2:B:88:LEU:HD21	3:B:147:HEM:CBA	2.19	0.45
2:B:93:CYS:HB3	2:B:94:ASP:OD2	2.16	0.45
2:B:64:GLY:O	2:B:68:LEU:HB2	2.17	0.45
1:A:89:HIS:CG	1:A:139:LYS:HE2	2.52	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:86:LEU:O	1:A:91:LEU:HG	2.17	0.44
1:A:37:PRO:O	1:A:40:LYS:HG3	2.18	0.44
1:A:109:LEU:O	1:A:113:LEU:HB2	2.17	0.44
2:B:63:HIS:CE1	3:B:147:HEM:CAD	3.00	0.44
1:A:11:LYS:HE3	1:A:73:LEU:HD12	2.00	0.44
1:A:53:ALA:HA	1:A:56:LYS:HB2	2.00	0.44
2:B:67:VAL:C	2:B:69:HIS:H	2.21	0.44
2:B:4:SER:HB3	2:B:7:GLU:CG	2.48	0.44
2:B:92:HIS:HE1	3:B:147:HEM:C4A	2.35	0.44
1:A:27:GLU:O	1:A:31:ARG:HG3	2.19	0.43
1:A:95:PRO:HA	1:A:98:PHE:CD2	2.52	0.43
2:B:98:VAL:HG11	2:B:103:PHE:HE1	1.83	0.43
1:A:7:LYS:O	1:A:11:LYS:HG3	2.19	0.43
2:B:51:PRO:O	2:B:55:MET:HG2	2.19	0.43
2:B:127:GLN:O	2:B:131:GLN:HG2	2.18	0.43
1:A:47:ASP:N	1:A:54:GLN:OE1	2.40	0.43
2:B:62:ALA:HB3	2:B:63:HIS:H	1.40	0.43
2:B:127:GLN:O	2:B:128:ALA:C	2.55	0.43
2:B:123:THR:O	2:B:127:GLN:N	2.34	0.42
2:B:128:ALA:O	2:B:131:GLN:HB2	2.18	0.42
2:B:67:VAL:C	2:B:69:HIS:N	2.70	0.42
1:A:89:HIS:HB2	1:A:139:LYS:HE2	2.01	0.42
1:A:94:ASP:HA	1:A:95:PRO:HD2	1.86	0.42
2:B:137:VAL:C	2:B:139:ASN:N	2.72	0.42
2:B:56:GLY:O	2:B:57:ASN:C	2.55	0.42
2:B:102:ASN:O	2:B:103:PHE:C	2.56	0.42
2:B:45:PHE:N	2:B:45:PHE:CD2	2.87	0.42
2:B:123:THR:OG1	2:B:126:LEU:HB2	2.19	0.42
1:A:36:PHE:CD2	1:A:100:LEU:HD22	2.55	0.42
2:B:93:CYS:HB2	2:B:145:TYR:CD2	2.55	0.42
1:A:31:ARG:NH2	2:B:127:GLN:OE1	2.52	0.42
2:B:103:PHE:N	2:B:103:PHE:CD1	2.87	0.42
1:A:83:LEU:HD21	3:A:142:HEM:HMA1	2.02	0.42
2:B:95:LYS:NZ	2:B:95:LYS:HB3	2.33	0.42
2:B:3:LEU:HD12	2:B:8:LYS:HE3	2.02	0.41
1:A:54:GLN:O	1:A:58:HIS:N	2.53	0.41
1:A:114:PRO:O	2:B:116:ARG:NH2	2.49	0.41
2:B:29:GLY:O	2:B:33:VAL:HG23	2.19	0.41
2:B:42:PHE:CD2	3:B:147:HEM:HMD1	2.56	0.41
2:B:137:VAL:C	2:B:139:ASN:H	2.23	0.41
1:A:36:PHE:O	1:A:39:THR:HG23	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:65:LYS:O	2:B:66:LYS:O	2.39	0.41
1:A:36:PHE:CE2	1:A:100:LEU:HD22	2.55	0.41
1:A:113:LEU:N	1:A:114:PRO:CD	2.81	0.41
1:A:95:PRO:HA	1:A:98:PHE:HD2	1.86	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	139/141 (99%)	117 (84%)	18 (13%)	4 (3%)	4 24
2	B	144/146 (99%)	114 (79%)	27 (19%)	3 (2%)	7 33
All	All	283/287 (99%)	231 (82%)	45 (16%)	7 (2%)	5 28

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	4	SER
1	A	25	GLY
1	A	38	THR
1	A	41	THR
1	A	104	CYS
2	B	140	ALA
2	B	64	GLY

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	115/115 (100%)	89 (77%)	26 (23%)	1 4
2	B	118/118 (100%)	94 (80%)	24 (20%)	1 6
All	All	233/233 (100%)	183 (78%)	50 (22%)	1 5

All (50) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1	VAL
1	A	7	LYS
1	A	9	ASN
1	A	16	LYS
1	A	23	GLU
1	A	34	LEU
1	A	38	THR
1	A	48	LEU
1	A	49	SER
1	A	56	LYS
1	A	60	LYS
1	A	61	LYS
1	A	81	SER
1	A	85	ASP
1	A	89	HIS
1	A	92	ARG
1	A	96	VAL
1	A	101	LEU
1	A	102	SER
1	A	105	LEU
1	A	109	LEU
1	A	113	LEU
1	A	125	LEU
1	A	136	LEU
1	A	138	SER
1	A	139	LYS
2	B	1	VAL
2	B	2	GLN
2	B	6	GLU
2	B	8	LYS
2	B	20	GLU
2	B	21	GLU
2	B	28	LEU

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Mol	Chain	Res	Type
2	B	30	ARG
2	B	61	LYS
2	B	65	LYS
2	B	68	LEU
2	B	71	PHE
2	B	73	GLU
2	B	81	LEU
2	B	82	LYS
2	B	94	ASP
2	B	95	LYS
2	B	104	ARG
2	B	124	PRO
2	B	125	GLU
2	B	126	LEU
2	B	144	LYS
2	B	145	TYR
2	B	146	HIS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	122	HIS
2	B	63	HIS
2	B	80	ASN
2	B	131	GLN
2	B	143	HIS

5.3.3 RNA (i)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates (i)

There are no monosaccharides in this entry.

5.6 Ligand geometry (i)

2 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	$\#\|Z\| > 2$
3	HEM	A	142	1	41,50,50	2.59	18 (43%)	45,82,82	2.00	10 (22%)
3	HEM	B	147	2	41,50,50	3.03	21 (51%)	45,82,82	2.92	21 (46%)

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
3	HEM	A	142	1	-	7/12/54/54	-
3	HEM	B	147	2	-	6/12/54/54	-

All (39) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	147	HEM	C3D-C2D	8.94	1.55	1.36
3	B	147	HEM	C1D-ND	6.57	1.52	1.38
3	A	142	HEM	C3C-C2C	-6.35	1.31	1.40
3	B	147	HEM	CAA-C2A	5.40	1.60	1.52
3	A	142	HEM	C3D-C2D	5.31	1.48	1.36
3	B	147	HEM	CMB-C2B	5.21	1.61	1.50
3	B	147	HEM	C4D-ND	5.02	1.49	1.40
3	B	147	HEM	FE-ND	4.55	2.19	1.96
3	B	147	HEM	C3C-CAC	4.29	1.56	1.47
3	A	142	HEM	CBD-CGD	4.29	1.60	1.50
3	A	142	HEM	CAA-C2A	4.25	1.58	1.52
3	A	142	HEM	C1B-NB	-4.16	1.33	1.40
3	A	142	HEM	C4A-CHB	-4.04	1.29	1.41
3	B	147	HEM	O1D-CGD	3.86	1.34	1.22
3	A	142	HEM	C3C-CAC	3.83	1.55	1.47

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	147	HEM	C4D-C3D	3.71	1.51	1.45
3	A	142	HEM	C3B-C2B	-3.66	1.29	1.37
3	A	142	HEM	FE-NB	-3.54	1.79	1.96
3	A	142	HEM	CAB-C3B	3.51	1.57	1.47
3	A	142	HEM	CMA-C3A	3.36	1.58	1.51
3	B	147	HEM	O1A-CGA	3.16	1.32	1.22
3	B	147	HEM	CAD-C3D	3.08	1.59	1.51
3	B	147	HEM	FE-NB	2.93	2.11	1.96
3	B	147	HEM	CAB-C3B	2.91	1.55	1.47
3	A	142	HEM	O1D-CGD	2.90	1.31	1.22
3	B	147	HEM	CMA-C3A	2.87	1.57	1.51
3	A	142	HEM	O1A-CGA	2.79	1.31	1.22
3	B	147	HEM	C1B-C2B	2.58	1.49	1.44
3	A	142	HEM	C4A-NA	-2.54	1.30	1.36
3	B	147	HEM	C1A-NA	2.41	1.41	1.36
3	A	142	HEM	CHB-C1B	-2.37	1.28	1.35
3	B	147	HEM	C4A-NA	2.36	1.41	1.36
3	A	142	HEM	C2A-C3A	-2.33	1.30	1.37
3	A	142	HEM	CMC-C2C	2.30	1.57	1.51
3	B	147	HEM	CMC-C2C	2.27	1.57	1.51
3	A	142	HEM	FE-ND	2.22	2.07	1.96
3	B	147	HEM	CBD-CGD	2.14	1.55	1.50
3	B	147	HEM	CBA-CGA	2.10	1.55	1.50
3	B	147	HEM	C3B-C4B	2.10	1.49	1.44

All (31) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
3	B	147	HEM	C2C-C3C-C4C	8.87	113.09	106.90
3	B	147	HEM	CHA-C4D-ND	-6.09	116.86	124.38
3	B	147	HEM	C3C-C4C-NC	-6.06	99.50	110.94
3	A	142	HEM	C2B-C1B-NB	5.97	116.91	109.84
3	A	142	HEM	C1B-NB-C4B	-5.57	99.32	105.07
3	B	147	HEM	CAD-C3D-C4D	4.96	133.32	124.66
3	B	147	HEM	CMC-C2C-C3C	4.59	133.27	124.68
3	B	147	HEM	CBD-CAD-C3D	-4.48	100.17	112.63
3	B	147	HEM	CMA-C3A-C4A	-4.30	121.85	128.46
3	B	147	HEM	C4C-CHD-C1D	3.64	127.36	122.56
3	B	147	HEM	CHA-C4D-C3D	3.64	132.16	125.33
3	A	142	HEM	CAD-CBD-CGD	3.57	121.30	113.60
3	B	147	HEM	O2D-CGD-O1D	3.48	131.98	123.30
3	B	147	HEM	O1D-CGD-CBD	-3.43	112.06	123.08

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
3	A	142	HEM	CBD-CAD-C3D	-3.29	103.48	112.63
3	B	147	HEM	CMA-C3A-C2A	3.03	130.65	124.94
3	B	147	HEM	C1D-C2D-C3D	3.01	110.11	106.96
3	B	147	HEM	CBA-CAA-C2A	-2.99	107.52	112.62
3	B	147	HEM	O1A-CGA-CBA	-2.90	113.76	123.08
3	A	142	HEM	CAA-C2A-C3A	-2.89	118.95	127.25
3	A	142	HEM	C4B-C3B-C2B	-2.67	105.00	107.11
3	A	142	HEM	CMD-C2D-C3D	2.56	133.07	126.12
3	B	147	HEM	CAD-C3D-C2D	-2.53	123.16	127.88
3	B	147	HEM	O2A-CGA-O1A	2.44	129.39	123.30
3	B	147	HEM	C3B-C2B-C1B	-2.35	104.74	106.49
3	B	147	HEM	C4D-C3D-C2D	-2.34	103.49	106.90
3	A	142	HEM	CMD-C2D-C1D	-2.20	121.69	125.04
3	B	147	HEM	CHD-C1D-C2D	2.16	128.35	124.98
3	A	142	HEM	CMA-C3A-C2A	2.12	128.94	124.94
3	A	142	HEM	CHB-C1B-C2B	-2.09	120.94	126.72
3	B	147	HEM	C2D-C1D-ND	-2.01	107.47	109.88

There are no chirality outliers.

All (13) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	142	HEM	C1A-C2A-CAA-CBA
3	A	142	HEM	C3A-C2A-CAA-CBA
3	A	142	HEM	C2B-C3B-CAB-CBB
3	A	142	HEM	C4B-C3B-CAB-CBB
3	B	147	HEM	C2B-C3B-CAB-CBB
3	B	147	HEM	CAA-CBA-CGA-O2A
3	B	147	HEM	CAA-CBA-CGA-O1A
3	B	147	HEM	C3D-CAD-CBD-CGD
3	A	142	HEM	CAA-CBA-CGA-O2A
3	B	147	HEM	C4B-C3B-CAB-CBB
3	A	142	HEM	CAA-CBA-CGA-O1A
3	A	142	HEM	C4D-C3D-CAD-CBD
3	B	147	HEM	CAD-CBD-CGD-O2D

There are no ring outliers.

2 monomers are involved in 30 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	142	HEM	3	0

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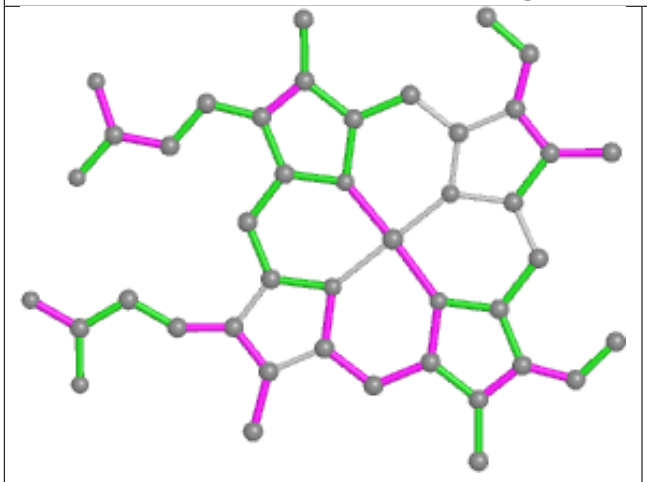
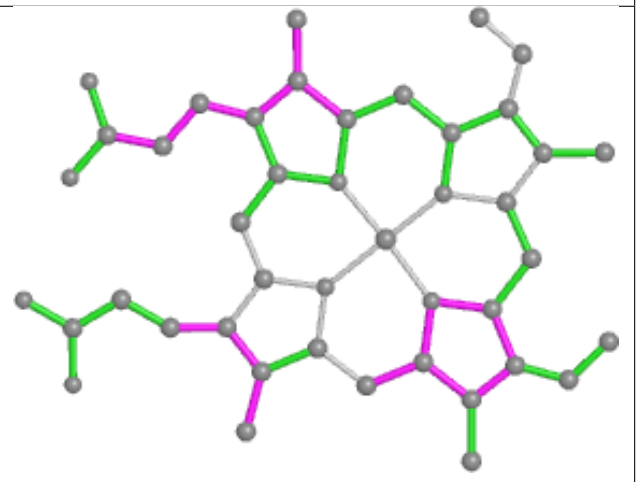
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	147	HEM	27	0

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 then 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.

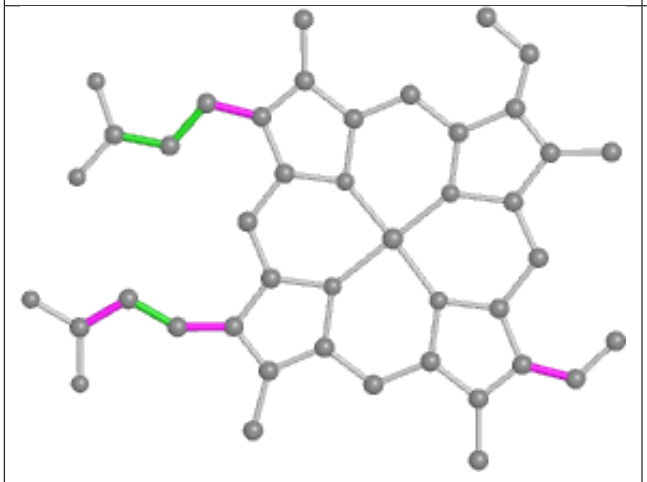
Ligand HEM A 142



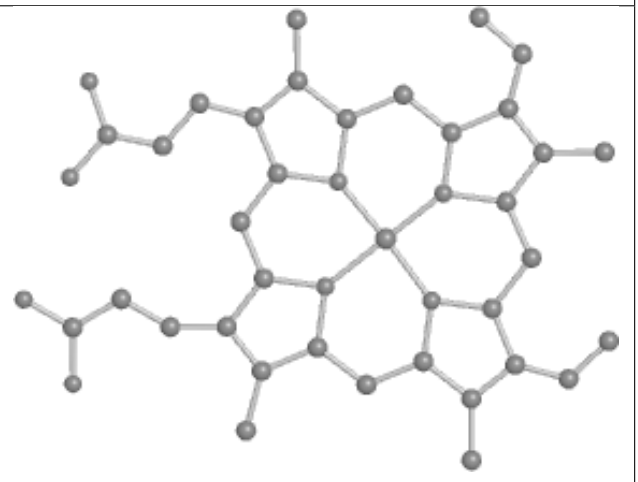
Bond lengths



Bond angles

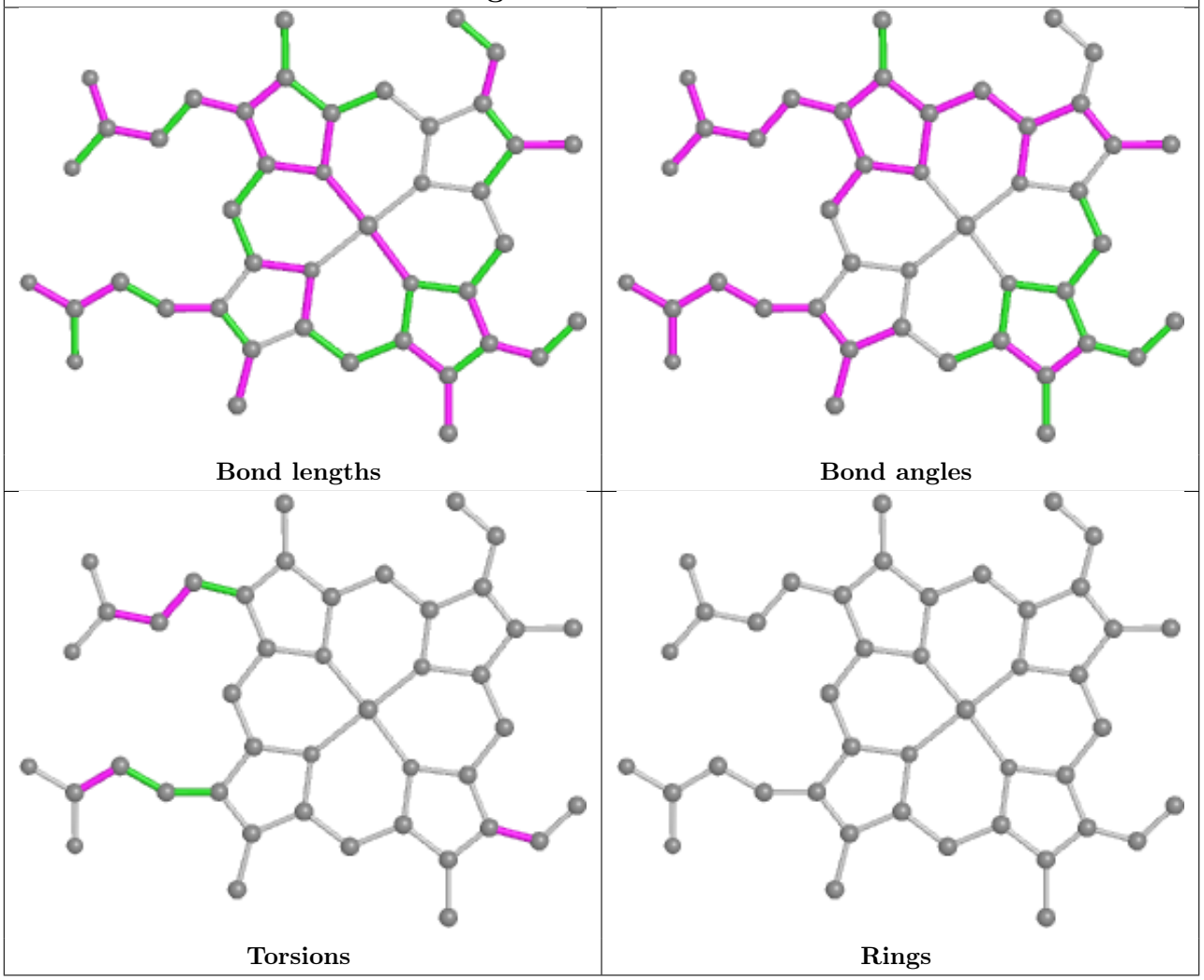


Torsions



Rings

Ligand HEM B 147



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	A	2

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	A	58:HIS	C	59:GLY	N	1.10

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Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	A	57:ALA	C	58:HIS	N	1.09

6 Fit of model and data

6.1 Protein, DNA and RNA chains

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, 95^{th} 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	141/141 (100%)	-0.76	1 (0%)	87	69	14, 29, 50, 80	0
2	B	146/146 (100%)	-0.48	2 (1%)	75	49	11, 35, 61, 89	1 (0%)
All	All	287/287 (100%)	-0.62	3 (1%)	82	59	11, 32, 59, 89	1 (0%)

All (3) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	146	HIS	3.0
1	A	141	ARG	2.3
2	B	94	ASP	2.2

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

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

6.3 Carbohydrates

There are no monosaccharides in this entry.

6.4 Ligands

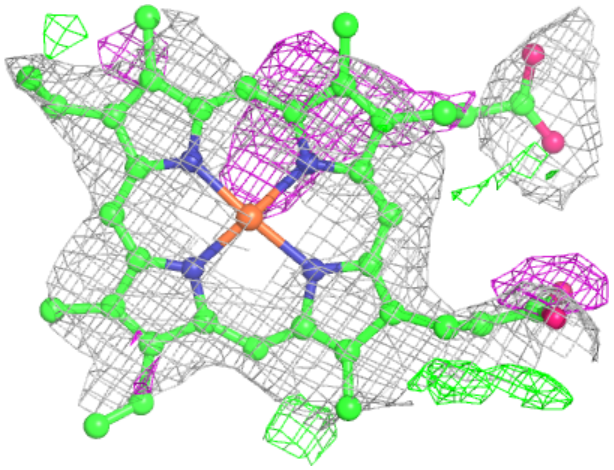
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95^{th} percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9'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

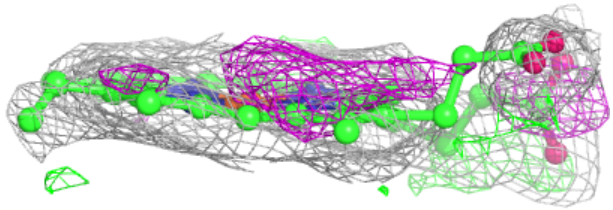
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
3	HEM	B	147	43/43	0.85	0.25	20,31,43,50	0
3	HEM	A	142	43/43	0.96	0.16	15,30,53,64	0

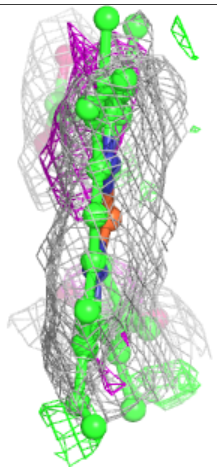
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around HEM B 147:

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

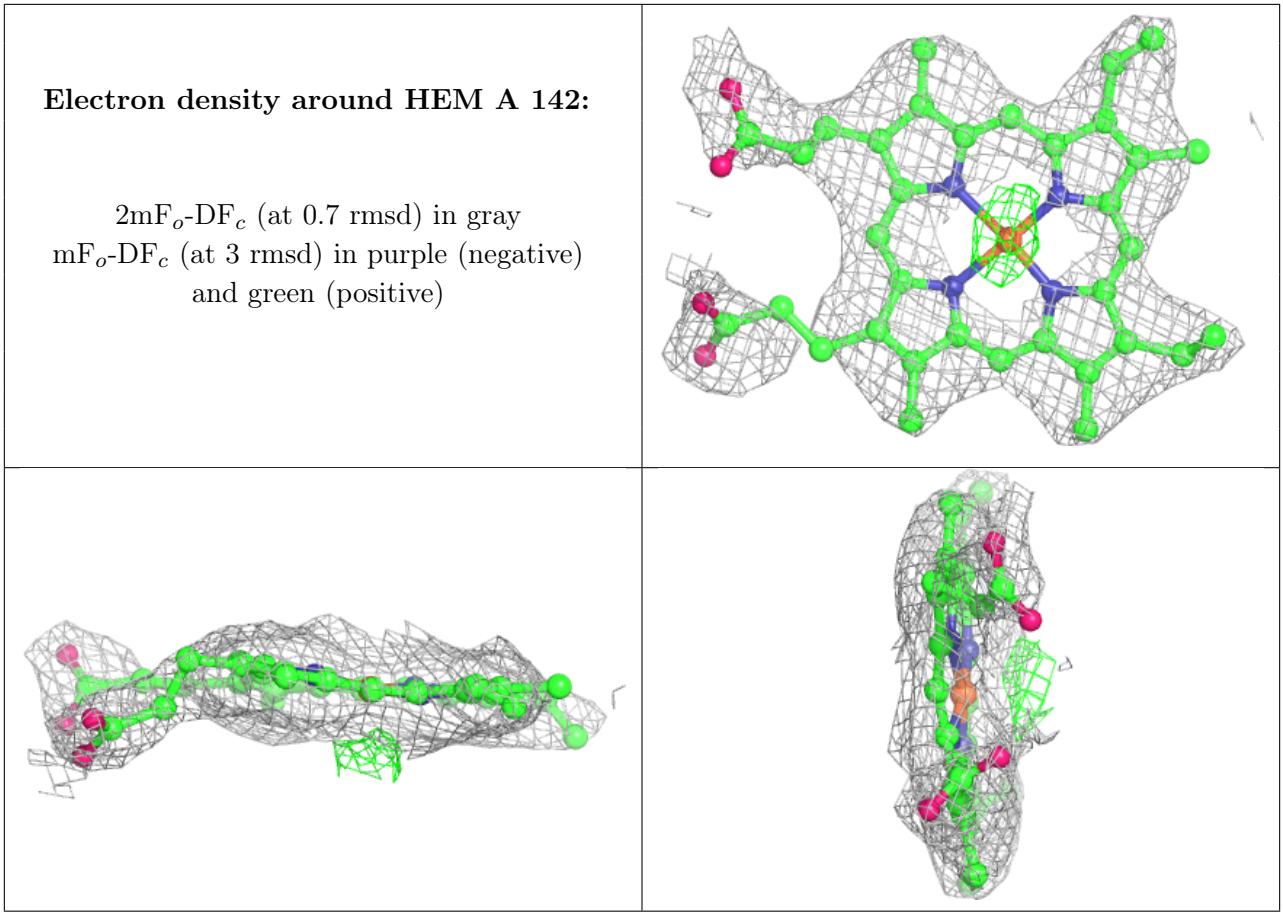






Electron density around HEM A 142:

$2mF_o$-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

There are no such residues in this entry.