



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 08:28 PM UTC

PDB ID : 2VV8 / pdb_00002vv8
Title : Co-bound structure of bjFixLH
Authors : Ayers, R.A.; Moffat, K.
Deposited on : 2008-06-04
Resolution : 1.61 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

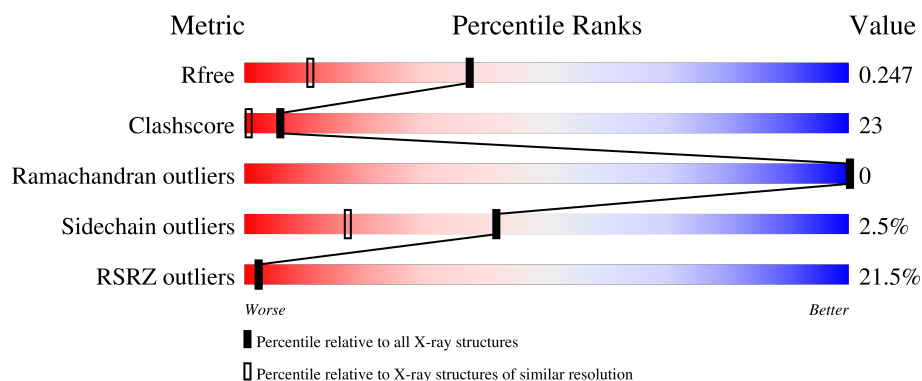
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.61 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	6728 (1.64-1.60)
Clashscore	190562	7023 (1.64-1.60)
Ramachandran outliers	187476	6898 (1.64-1.60)
Sidechain outliers	187428	6896 (1.64-1.60)
RSRZ outliers	180081	6727 (1.64-1.60)

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	119	19% 63% 24% • 11%
1	B	119	17% 65% 24% • 10%
1	C	119	18% 67% 21% • 11%
1	D	119	23% 71% 18% • 9%

2 Entry composition [i](#)

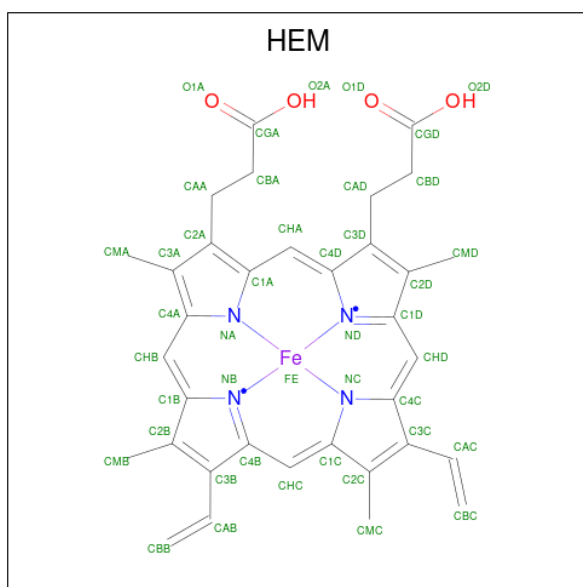
There are 6 unique types of molecules in this entry. The entry contains 4190 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 SENSOR PROTEIN FIXL.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	106	Total	C	N	O	S	0	15	0
			925	595	160	163	7			
1	B	107	Total	C	N	O	S	0	9	1
			882	559	158	159	6			
1	C	106	Total	C	N	O	S	0	6	1
			859	544	154	156	5			
1	D	108	Total	C	N	O	S	0	6	1
			873	554	153	161	5			

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



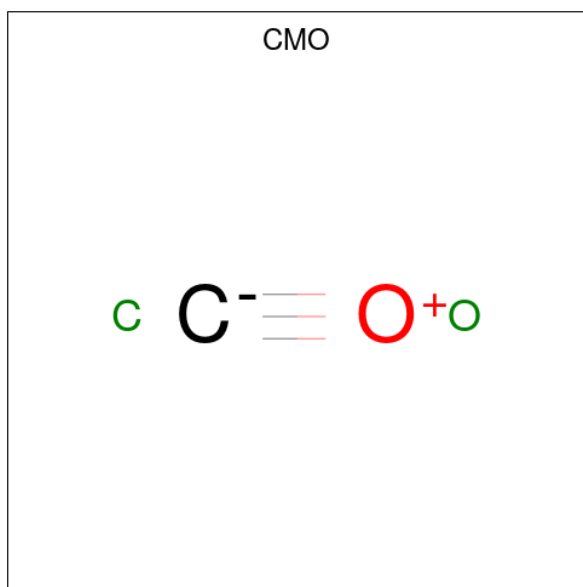
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	Fe	N O	0	0
			43	34	1	4 4		
2	B	1	Total	C	Fe	N O	0	0
			43	34	1	4 4		

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

- Molecule 3 is CARBON MONOXIDE (CCD ID: CMO) (formula: CO).



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

- Molecule 4 is SODIUM ION (CCD ID: NA) (formula: Na).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	C	1	Total	Na		
			1	1		

- Molecule 5 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	C	2	Total	Cl	0	0
			2	2		

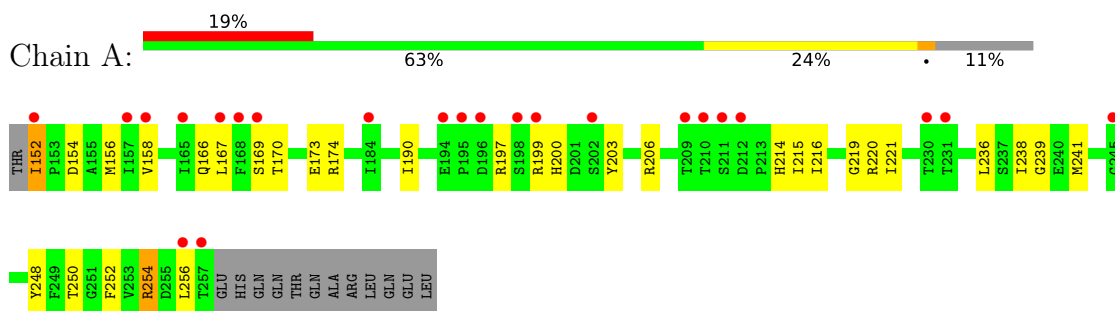
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	126	Total	O	0	0
			126	126		
6	B	120	Total	O	0	0
			120	120		
6	C	125	Total	O	0	0
			125	125		
6	D	97	Total	O	0	0
			97	97		

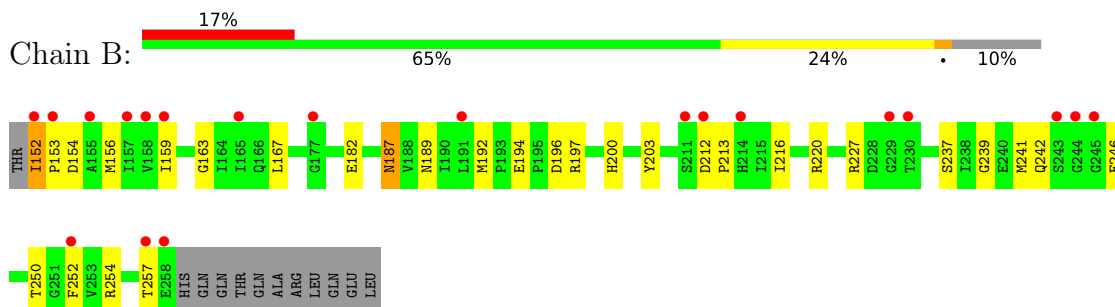
3 Residue-property plots [i](#)

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.

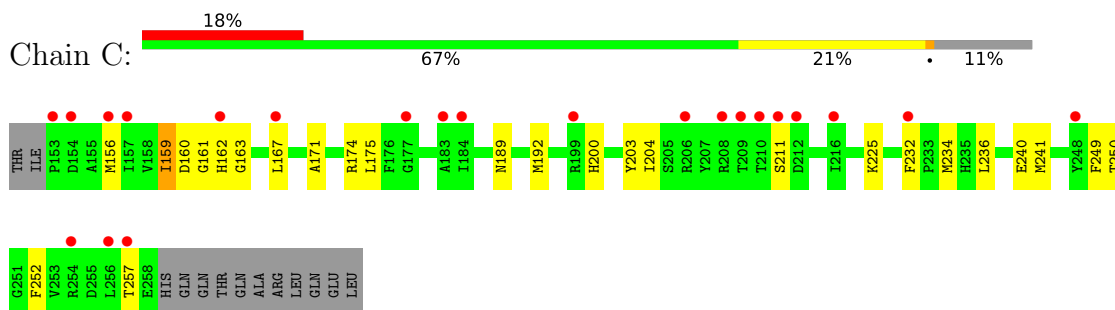
• Molecule 1: SENSOR PROTEIN FIXL



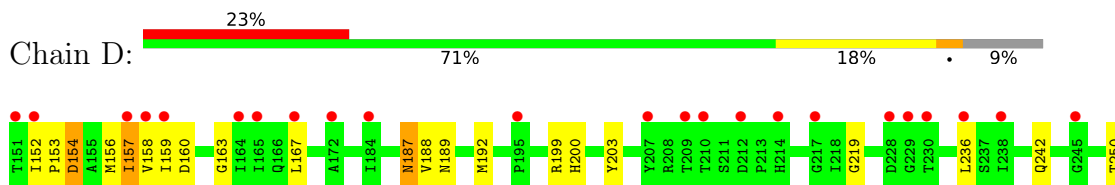
• Molecule 1: SENSOR PROTEIN FIXL



• Molecule 1: SENSOR PROTEIN FIXL



• Molecule 1: SENSOR PROTEIN FIXL



G251	F252	V253	R254	D255	L256	T257	E258	HIS	GLN	GLN	THR	GLN	ALA	ARG	LEU	GLN	GLU	LEU
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4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	48.92Å 50.30Å 58.99Å 73.37° 70.26° 72.13°	Depositor
Resolution (Å)	46.88 – 1.61 46.88 – 1.61	Depositor EDS
% Data completeness (in resolution range)	91.6 (46.88-1.61) 81.1 (46.88-1.61)	Depositor EDS
R_{merge}	0.04	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.55 (at 1.61Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.206 , 0.249 0.205 , 0.247	Depositor DCC
R_{free} test set	2619 reflections (4.90%)	wwPDB-VP
Wilson B-factor (Å ²)	20.1	Xtriage
Anisotropy	0.031	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 43.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.45$, $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	4190	wwPDB-VP
Average B, all atoms (Å ²)	16.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.79% 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, CMO, 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	0.71	0/998	0.92	2/1345 (0.1%)
1	B	0.77	1/928 (0.1%)	0.89	0/1253
1	C	0.82	1/896 (0.1%)	1.06	2/1209 (0.2%)
1	D	0.85	1/910 (0.1%)	0.90	2/1232 (0.2%)
All	All	0.79	3/3732 (0.1%)	0.95	6/5039 (0.1%)

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	C	0	2

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	257	THR	C-N	-7.62	1.22	1.33
1	D	257	THR	C-N	-6.67	1.24	1.33
1	B	257	THR	C-N	-6.41	1.24	1.33

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	162	HIS	N-CA-C	13.84	128.20	112.57
1	C	160	ASP	N-CA-C	10.55	123.94	110.33
1	A	152	ILE	CA-C-N	-6.14	113.64	120.14
1	A	152	ILE	C-N-CA	-6.14	113.64	120.14
1	D	152	ILE	CA-C-N	5.01	125.28	119.92
1	D	152	ILE	C-N-CA	5.01	125.28	119.92

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	C	159	ILE	Peptide
1	C	161	GLY	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	925	0	979	72	0
1	B	882	0	899	57	0
1	C	859	0	866	32	0
1	D	873	0	878	35	0
2	A	43	0	30	5	0
2	B	43	0	30	4	0
2	C	43	0	30	2	0
2	D	43	0	30	1	0
3	A	2	0	0	0	0
3	B	2	0	0	0	0
3	C	2	0	0	0	0
3	D	2	0	0	0	0
4	C	1	0	0	0	0
5	C	2	0	0	0	0
6	A	126	0	0	13	0
6	B	120	0	0	3	0
6	C	125	0	0	6	0
6	D	97	0	0	3	0
All	All	4190	0	3742	171	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 (171) 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:D:159[B]:ILE:CD1	1:D:188:VAL:HB	1.74	1.17

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:156[A]:MET:HE3	1:B:156:MET:HE3	1.27	1.16
1:D:159[B]:ILE:HD11	1:D:188:VAL:HB	1.23	1.13
1:A:158[B]:VAL:CG2	1:A:167[B]:LEU:HB3	1.78	1.12
1:A:170:THR:O	1:A:173[B]:GLU:HG2	1.50	1.12
1:A:156[A]:MET:CE	1:B:156:MET:HE3	1.79	1.10
1:C:156:MET:HE1	1:D:158:VAL:HG21	1.25	1.09
1:A:174[A]:ARG:HD3	6:A:2029:HOH:O	1.53	1.07
1:C:175:LEU:HD11	1:C:232:PHE:CZ	1.90	1.06
1:A:158[B]:VAL:HG22	1:A:167[B]:LEU:HB3	1.06	1.05
1:C:156:MET:HE2	1:C:252:PHE:CE1	1.95	1.01
1:A:156[A]:MET:HE3	1:B:156:MET:CE	1.89	1.01
1:A:252:PHE:HE1	1:B:241[A]:MET:SD	1.84	1.01
1:B:153:PRO:HG3	1:B:254[B]:ARG:HE	1.29	0.94
1:B:152:ILE:HA	6:B:2001:HOH:O	1.68	0.92
1:A:239:GLY:HA3	1:B:241[B]:MET:HG3	1.51	0.91
1:B:187:ASN:HD22	1:B:189:ASN:H	1.18	0.88
1:A:239:GLY:CA	1:B:241[B]:MET:HG3	2.02	0.87
1:A:167[A]:LEU:HD22	1:B:156:MET:HE1	1.55	0.86
1:A:170:THR:HG22	6:A:2022:HOH:O	1.74	0.85
1:C:156:MET:HE3	1:D:167:LEU:HD13	1.55	0.85
1:A:254[A]:ARG:HG2	1:A:254[A]:ARG:HH21	1.42	0.84
1:D:187:ASN:HD22	1:D:189:ASN:H	1.25	0.84
1:D:154:ASP:HB3	6:D:2007:HOH:O	1.78	0.84
1:D:159[B]:ILE:HD11	1:D:188:VAL:CB	2.05	0.84
1:B:241[A]:MET:HE3	1:B:250[A]:THR:HG23	1.58	0.83
1:D:153:PRO:HG3	1:D:254[B]:ARG:HG3	1.60	0.83
1:A:252:PHE:CE1	1:B:241[A]:MET:SD	2.72	0.82
1:B:241[A]:MET:CE	1:B:250[A]:THR:CG2	2.57	0.82
1:B:241[A]:MET:CE	1:B:250[A]:THR:HG21	2.10	0.81
1:A:156[B]:MET:HG3	1:A:252:PHE:CE2	2.16	0.81
1:D:159[B]:ILE:HG21	1:D:163:GLY:HA2	1.62	0.81
1:A:156[C]:MET:SD	1:B:156:MET:HE1	2.21	0.80
1:C:156:MET:HE2	1:C:252:PHE:HE1	1.49	0.78
1:D:156:MET:C	1:D:157[A]:ILE:HD12	2.08	0.78
1:B:213:PRO:HB3	1:B:216[B]:ILE:HB	1.65	0.77
1:A:250[B]:THR:HG23	6:A:2122:HOH:O	1.84	0.75
6:C:2001:HOH:O	1:D:156:MET:HE1	1.86	0.75
1:A:174[A]:ARG:CD	6:A:2029:HOH:O	2.22	0.74
1:A:173[B]:GLU:CD	1:A:174[B]:ARG:NH2	2.47	0.73
1:B:241[A]:MET:HE3	1:B:250[A]:THR:CG2	2.17	0.73
1:C:156:MET:CE	1:D:158:VAL:HG21	2.13	0.72

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:216:ILE:HG13	1:A:238[B]:ILE:HD12	1.70	0.72
1:A:254[A]:ARG:HH21	1:A:254[A]:ARG:CG	2.03	0.71
1:B:237[B]:SER:OG	1:B:252:PHE:HB2	1.91	0.70
1:A:254[A]:ARG:HG2	1:A:254[A]:ARG:NH2	2.05	0.70
1:B:241[A]:MET:HE1	1:B:250[A]:THR:HG21	1.72	0.70
1:B:192[B]:MET:HE2	1:B:196:ASP:HB3	1.74	0.69
1:C:175:LEU:CD1	1:C:232:PHE:CZ	2.73	0.69
1:A:216:ILE:HA	1:A:238[B]:ILE:HG13	1.74	0.68
1:A:254[A]:ARG:HD2	6:A:2003:HOH:O	1.93	0.68
1:D:159[B]:ILE:CD1	1:D:188:VAL:CB	2.64	0.66
1:C:156:MET:CE	1:D:167:LEU:HD13	2.24	0.66
1:C:175:LEU:HD11	1:C:232:PHE:CE1	2.30	0.66
1:A:173[B]:GLU:OE2	1:A:174[B]:ARG:NH2	2.29	0.66
1:A:154:ASP:OD2	1:A:254[A]:ARG:HG3	1.95	0.66
1:A:241[B]:MET:HB3	1:B:239:GLY:HA3	1.77	0.65
1:A:158[B]:VAL:HG12	1:A:250[B]:THR:HG22	1.78	0.63
1:B:241[A]:MET:CE	1:B:250[A]:THR:HG23	2.23	0.63
1:D:199:ARG:NH1	6:D:2043:HOH:O	2.29	0.63
1:C:250:THR:HG23	6:C:2117:HOH:O	1.98	0.63
1:C:175:LEU:C	1:C:175:LEU:HD13	2.25	0.62
1:B:154:ASP:HB2	6:B:2007:HOH:O	2.00	0.62
1:B:152:ILE:N	1:B:153:PRO:HD3	2.15	0.61
1:A:206[A]:ARG:HH22	1:A:214:HIS:CE1	2.18	0.61
1:A:250[B]:THR:HG21	1:B:250[B]:THR:HG21	1.81	0.61
1:D:159[B]:ILE:CG2	1:D:163:GLY:HA2	2.31	0.61
1:A:156[B]:MET:HE2	1:B:167:LEU:HD13	1.80	0.61
1:A:170:THR:HG21	6:A:2016:HOH:O	2.01	0.60
1:A:241[A]:MET:HE3	6:A:2115:HOH:O	1.99	0.60
1:B:246:GLU:HG3	6:B:2113:HOH:O	2.02	0.59
1:A:167[B]:LEU:CD1	1:B:156:MET:HE1	2.33	0.59
1:B:189:ASN:HA	1:B:192[A]:MET:HE3	1.85	0.58
1:D:159[B]:ILE:HG22	1:D:160:ASP:O	2.03	0.58
1:C:200:HIS:CE1	1:C:204[B]:ILE:HD11	2.39	0.58
1:A:221[B]:ILE:HD12	1:A:256:LEU:HD11	1.84	0.58
1:A:158[B]:VAL:CG2	1:A:167[B]:LEU:CB	2.71	0.57
1:A:166:GLN:NE2	6:A:2011:HOH:O	2.37	0.57
1:B:192[B]:MET:CE	1:B:196:ASP:HB3	2.35	0.57
1:B:187:ASN:ND2	1:B:189:ASN:H	1.96	0.56
1:C:159:ILE:HD12	1:C:163:GLY:HA2	1.87	0.56
1:C:171:ALA:HA	1:C:174[B]:ARG:HG2	1.87	0.56
1:C:241[B]:MET:SD	1:D:252:PHE:CE1	2.98	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:156[B]:MET:HG3	1:A:252:PHE:CD2	2.40	0.56
1:B:152:ILE:N	1:B:153:PRO:CD	2.69	0.56
1:A:215[A]:ILE:CD1	2:A:1258:HEM:HBA1	2.36	0.55
2:A:1258:HEM:HBC2	2:A:1258:HEM:HMC2	1.88	0.55
1:A:200:HIS:HA	1:A:203:TYR:CD2	2.41	0.55
2:D:1259:HEM:HMC1	2:D:1259:HEM:HBC2	1.88	0.55
1:B:213:PRO:HB3	1:B:216[A]:ILE:HB	1.88	0.54
1:B:189:ASN:HA	1:B:192[B]:MET:SD	2.48	0.54
1:B:194:GLU:OE2	1:B:197[B]:ARG:HD2	2.08	0.54
1:A:167[A]:LEU:HD22	1:B:156:MET:CE	2.35	0.54
1:A:206[A]:ARG:NH2	1:A:214:HIS:CE1	2.76	0.53
1:A:216:ILE:HG13	1:A:238[B]:ILE:CD1	2.36	0.53
1:C:175:LEU:HD11	1:C:232:PHE:HZ	1.65	0.53
1:A:190:ILE:HA	1:A:197:ARG:NH2	2.24	0.53
1:A:156[B]:MET:HB3	1:A:169[B]:SER:OG	2.09	0.53
1:A:170:THR:O	1:A:173[B]:GLU:CG	2.41	0.53
1:C:241[A]:MET:HG2	6:C:2117:HOH:O	2.08	0.52
1:B:241[B]:MET:HG2	1:B:242:GLN:H	1.75	0.52
6:A:2108:HOH:O	1:B:241[B]:MET:HG2	2.09	0.52
1:D:187:ASN:ND2	1:D:189:ASN:H	2.03	0.51
1:B:241[B]:MET:HG2	1:B:242:GLN:N	2.26	0.51
1:A:156[A]:MET:CE	1:B:156:MET:CE	2.63	0.51
1:D:159[B]:ILE:HG23	1:D:163:GLY:C	2.35	0.51
1:C:225:LYS:HD3	6:C:2039:HOH:O	2.10	0.50
1:A:250[B]:THR:CG2	6:A:2122:HOH:O	2.52	0.50
1:D:189:ASN:HA	1:D:192:MET:HE3	1.93	0.50
1:C:241[A]:MET:SD	1:D:252:PHE:HE1	2.35	0.49
1:C:234:MET:CE	1:C:236[B]:LEU:HD13	2.43	0.49
1:D:159[B]:ILE:CG2	1:D:163:GLY:CA	2.90	0.49
1:A:250[A]:THR:HB	1:B:241[A]:MET:HE1	1.94	0.48
1:C:200:HIS:O	1:C:204[B]:ILE:HG12	2.13	0.48
1:D:242:GLN:HG2	6:D:2089:HOH:O	2.13	0.48
1:C:156:MET:HE1	1:D:158:VAL:CG2	2.18	0.47
1:C:240:GLU:HG3	1:C:249:PHE:CE2	2.49	0.47
1:A:156[B]:MET:HE1	1:B:156:MET:HE3	1.95	0.47
2:B:1259:HEM:HBB2	2:B:1259:HEM:HHC	1.96	0.47
1:A:241[B]:MET:SD	1:A:248:TYR:HB2	2.54	0.47
1:C:189:ASN:HA	1:C:192:MET:HE3	1.96	0.47
1:A:156[B]:MET:CE	1:B:156:MET:HE3	2.44	0.47
1:A:156[A]:MET:HE1	1:B:156:MET:HE3	1.86	0.47
1:B:182:GLU:OE1	1:B:227:ARG:NH2	2.48	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:159[B]:ILE:HD11	1:D:188:VAL:CG2	2.44	0.46
1:C:234:MET:HE2	1:C:236[B]:LEU:HD13	1.96	0.46
1:B:154:ASP:OD1	1:B:154:ASP:N	2.46	0.46
1:A:215[A]:ILE:CD1	2:A:1258:HEM:CBA	2.94	0.46
1:C:167:LEU:HG	6:C:2010:HOH:O	2.16	0.45
1:D:250[B]:THR:HG22	1:D:252:PHE:CE1	2.51	0.45
1:A:167[A]:LEU:CD1	1:B:156:MET:HE2	2.46	0.45
1:A:158[B]:VAL:CG1	1:A:250[B]:THR:HG22	2.47	0.45
2:A:1258:HEM:HBC2	2:A:1258:HEM:CMC	2.47	0.45
1:A:167[A]:LEU:HD13	1:B:156:MET:HE2	1.99	0.45
1:B:200:HIS:HA	1:B:203:TYR:CD1	2.52	0.44
2:B:1259:HEM:HBC2	2:B:1259:HEM:HMC2	2.00	0.44
1:C:200:HIS:HA	1:C:203:TYR:CD2	2.53	0.44
1:C:241[A]:MET:SD	1:D:252:PHE:CE1	3.11	0.43
1:C:234:MET:CE	1:C:236[A]:LEU:HG	2.49	0.43
1:A:173[B]:GLU:CD	1:A:174[B]:ARG:HH21	2.24	0.43
1:B:153:PRO:HG3	1:B:254[B]:ARG:NE	2.14	0.43
1:D:158:VAL:HG22	1:D:250[A]:THR:HG22	2.01	0.43
1:A:215[A]:ILE:HD11	1:A:220:ARG:CZ	2.49	0.43
1:A:167[A]:LEU:CD2	1:B:156:MET:HE1	2.37	0.42
1:B:192[B]:MET:HE3	2:B:1259:HEM:C2B	2.53	0.42
1:A:254[A]:ARG:HD3	6:A:2123:HOH:O	2.18	0.42
1:B:159:ILE:HD12	1:B:163:GLY:HA2	1.99	0.42
1:A:206[C]:ARG:HD2	1:A:206[C]:ARG:HA	1.77	0.42
1:D:156:MET:HE3	1:D:167:LEU:CD2	2.50	0.42
1:B:192[B]:MET:HE1	1:B:200:HIS:HB2	2.00	0.42
1:A:199:ARG:HD3	6:A:2062:HOH:O	2.20	0.42
1:C:203:TYR:CD2	2:C:1259:HEM:C2A	3.07	0.42
1:A:156[A]:MET:HE3	1:B:156:MET:HE1	1.90	0.42
1:D:156:MET:HE3	1:D:167:LEU:HD21	2.03	0.41
1:A:241[B]:MET:HG3	6:A:2122:HOH:O	2.21	0.41
1:C:236[A]:LEU:HD11	2:C:1259:HEM:C4B	2.55	0.41
1:D:159[B]:ILE:HD12	1:D:188:VAL:HB	1.88	0.41
1:A:156[B]:MET:HB3	1:A:156[B]:MET:HE3	1.94	0.41
1:B:241[B]:MET:HE2	1:B:241[B]:MET:HB3	1.82	0.41
1:D:200:HIS:HA	1:D:203:TYR:CD1	2.55	0.41
1:A:219:GLY:HA2	1:A:236:LEU:O	2.20	0.41
2:A:1258:HEM:HHC	2:A:1258:HEM:HBB2	2.03	0.41
1:B:220:ARG:NH1	2:B:1259:HEM:HBA2	2.36	0.41
1:D:219:GLY:HA2	1:D:236:LEU:O	2.21	0.41
1:A:214:HIS:C	1:A:215[A]:ILE:HD12	2.45	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:241[B]:MET:HG2	6:C:2117:HOH:O	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	122/119 (102%)	120 (98%)	2 (2%)	0	100	100
1	B	114/119 (96%)	111 (97%)	3 (3%)	0	100	100
1	C	110/119 (92%)	108 (98%)	2 (2%)	0	100	100
1	D	112/119 (94%)	112 (100%)	0	0	100	100
All	All	458/476 (96%)	451 (98%)	7 (2%)	0	100	100

There are no Ramachandran outliers to report.

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	108/102 (106%)	105 (97%)	3 (3%)	38	14
1	B	99/102 (97%)	96 (97%)	3 (3%)	36	12
1	C	95/102 (93%)	94 (99%)	1 (1%)	65	46
1	D	97/102 (95%)	93 (96%)	4 (4%)	27	7

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
All	All	399/408 (98%)	388 (97%)	11 (3%)	42 14

All (11) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	152	ILE
1	A	254[A]	ARG
1	A	254[B]	ARG
1	B	152	ILE
1	B	187	ASN
1	B	212	ASP
1	C	211	SER
1	D	154	ASP
1	D	157[A]	ILE
1	D	157[B]	ILE
1	D	187	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	166	GLN
1	A	214	HIS
1	B	187	ASN
1	C	186	GLN
1	C	214	HIS
1	D	187	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

Of 11 ligands modelled in this entry, 3 are monoatomic - leaving 8 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	1259	1	50,50,50	1.81	9 (18%)	67,82,82	1.79	13 (19%)
3	CMO	A	1259	-	0,1,1	-	-	-	-	-
2	HEM	D	1259	1	50,50,50	1.91	10 (20%)	67,82,82	1.38	8 (11%)
3	CMO	C	1260	-	0,1,1	-	-	-	-	-
3	CMO	D	1260	-	0,1,1	-	-	-	-	-
3	CMO	B	1260	-	0,1,1	-	-	-	-	-
2	HEM	A	1258	1	50,50,50	1.92	9 (18%)	67,82,82	2.30	23 (34%)
2	HEM	C	1259	1	50,50,50	1.60	5 (10%)	67,82,82	1.56	12 (17%)

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	D	1259	1	-	1/14/54/54	-
2	HEM	A	1258	1	-	3/14/54/54	-
2	HEM	C	1259	1	-	2/14/54/54	-
2	HEM	B	1259	1	-	4/14/54/54	-

All (33) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	1258	HEM	C3D-C2D	7.74	1.53	1.36
2	D	1259	HEM	C3D-C2D	7.65	1.53	1.36
2	B	1259	HEM	C3D-C2D	6.64	1.51	1.36
2	C	1259	HEM	C3D-C2D	6.10	1.49	1.36
2	B	1259	HEM	FE-NA	4.45	2.09	1.95
2	B	1259	HEM	FE-ND	4.28	2.08	1.94
2	D	1259	HEM	FE-NA	4.14	2.08	1.95

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	1258	HEM	FE-NA	4.04	2.08	1.95
2	D	1259	HEM	FE-ND	3.98	2.07	1.94
2	A	1258	HEM	FE-ND	3.78	2.06	1.94
2	C	1259	HEM	FE-NA	3.54	2.06	1.95
2	A	1258	HEM	CAB-C3B	3.47	1.56	1.47
2	D	1259	HEM	CAC-C3C	3.42	1.56	1.47
2	B	1259	HEM	FE-NC	3.39	2.06	1.95
2	C	1259	HEM	CAB-C3B	3.34	1.56	1.47
2	D	1259	HEM	FE-NC	3.28	2.06	1.95
2	B	1259	HEM	CAB-C3B	3.17	1.55	1.47
2	D	1259	HEM	CMB-C2B	2.95	1.56	1.50
2	C	1259	HEM	CAC-C3C	2.89	1.55	1.47
2	D	1259	HEM	CAB-C3B	2.84	1.55	1.47
2	A	1258	HEM	FE-NC	2.83	2.04	1.95
2	C	1259	HEM	FE-NC	2.78	2.04	1.95
2	A	1258	HEM	CAC-C3C	2.61	1.54	1.47
2	A	1258	HEM	CMA-C3A	2.55	1.56	1.50
2	D	1259	HEM	CMC-C2C	2.52	1.55	1.50
2	B	1259	HEM	CAC-C3C	2.51	1.54	1.47
2	D	1259	HEM	CMA-C3A	2.50	1.55	1.50
2	B	1259	HEM	CMA-C3A	2.44	1.55	1.50
2	D	1259	HEM	CMD-C2D	2.43	1.55	1.50
2	B	1259	HEM	CMC-C2C	2.37	1.55	1.50
2	A	1258	HEM	C1D-ND	2.23	1.43	1.38
2	A	1258	HEM	CAA-C2A	2.21	1.57	1.51
2	B	1259	HEM	C3B-C2B	-2.12	1.32	1.37

All (56) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	1258	HEM	C3B-C4B-NB	6.72	114.29	109.47
2	B	1259	HEM	CHD-C4C-NC	6.01	131.00	124.45
2	A	1258	HEM	CHC-C4B-NB	-5.38	118.63	124.42
2	B	1259	HEM	C3B-C2B-C1B	4.89	110.08	106.41
2	A	1258	HEM	CHD-C1D-ND	4.74	129.52	124.42
2	A	1258	HEM	C2A-C1A-NA	-4.73	104.90	110.15
2	A	1258	HEM	CAD-C3D-C4D	4.71	132.90	124.70
2	A	1258	HEM	C3B-C2B-C1B	4.51	109.80	106.41
2	A	1258	HEM	C4B-C3B-C2B	-4.17	103.45	107.28
2	A	1258	HEM	CHC-C1C-NC	4.14	128.96	124.45
2	D	1259	HEM	C4D-ND-C1D	4.13	110.10	105.21
2	C	1259	HEM	C3B-C2B-C1B	3.98	109.40	106.41

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	1259	HEM	CHA-C4D-ND	3.87	129.15	124.37
2	B	1259	HEM	C4C-NC-C1C	3.80	112.02	105.82
2	A	1258	HEM	C4A-NA-C1A	3.73	111.91	105.82
2	A	1258	HEM	C4A-C3A-C2A	3.61	110.95	106.82
2	C	1259	HEM	C3B-C4B-NB	3.56	112.03	109.47
2	B	1259	HEM	C4C-C3C-C2C	3.53	109.88	106.81
2	B	1259	HEM	C4D-ND-C1D	3.52	109.38	105.21
2	C	1259	HEM	CAD-C3D-C4D	3.45	130.71	124.70
2	A	1258	HEM	CMA-C3A-C4A	-3.39	120.26	125.42
2	B	1259	HEM	CAA-CBA-CGA	-3.37	104.71	113.67
2	A	1258	HEM	CHD-C1D-C2D	-3.23	119.93	125.03
2	A	1258	HEM	CAD-C3D-C2D	-3.21	121.86	127.87
2	C	1259	HEM	C4D-ND-C1D	3.17	108.97	105.21
2	A	1258	HEM	C3A-C4A-NA	-3.03	105.28	110.14
2	B	1259	HEM	CMA-C3A-C4A	-2.92	120.98	125.42
2	A	1258	HEM	CHA-C1A-NA	2.85	129.03	123.86
2	D	1259	HEM	C3B-C2B-C1B	2.80	108.51	106.41
2	A	1258	HEM	CHA-C4D-ND	-2.79	120.91	124.37
2	C	1259	HEM	C4C-C3C-C2C	2.70	109.16	106.81
2	B	1259	HEM	CAC-C3C-C4C	-2.70	118.37	124.82
2	C	1259	HEM	CAA-CBA-CGA	-2.64	106.66	113.67
2	A	1258	HEM	CMD-C2D-C1D	2.49	128.93	125.03
2	C	1259	HEM	C4B-C3B-C2B	-2.44	105.03	107.28
2	A	1258	HEM	C4C-C3C-C2C	2.44	108.93	106.81
2	D	1259	HEM	CHA-C4D-ND	2.43	127.38	124.37
2	A	1258	HEM	C4D-ND-C1D	2.42	108.08	105.21
2	B	1259	HEM	O2A-CGA-CBA	2.38	121.53	114.00
2	D	1259	HEM	CBD-CAD-C3D	-2.38	105.95	112.53
2	D	1259	HEM	O2D-CGD-CBD	2.36	121.46	114.00
2	C	1259	HEM	CHD-C1D-ND	2.35	126.95	124.42
2	C	1259	HEM	CBD-CAD-C3D	-2.29	106.19	112.53
2	B	1259	HEM	CHB-C4A-NA	2.29	128.01	123.86
2	B	1259	HEM	CAD-C3D-C4D	2.29	128.69	124.70
2	C	1259	HEM	C3C-C2C-C1C	-2.27	104.90	107.05
2	A	1258	HEM	O1D-CGD-CBD	-2.27	115.91	123.09
2	D	1259	HEM	O1D-CGD-CBD	-2.22	116.06	123.09
2	B	1259	HEM	C4A-C3A-C2A	2.18	109.31	106.82
2	D	1259	HEM	C2A-C1A-NA	-2.18	107.74	110.15
2	A	1258	HEM	C1B-NB-C4B	-2.18	102.63	105.21
2	C	1259	HEM	CAC-C3C-C4C	-2.17	119.64	124.82
2	A	1258	HEM	CHD-C4C-C3C	2.14	128.81	125.21
2	D	1259	HEM	CHD-C1D-ND	2.04	126.62	124.42

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	1259	HEM	CHD-C4C-C3C	2.03	128.64	125.21
2	A	1258	HEM	O2D-CGD-CBD	2.03	120.41	114.00

There are no chirality outliers.

All (10) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	B	1259	HEM	C3D-CAD-CBD-CGD
2	D	1259	HEM	C3D-CAD-CBD-CGD
2	B	1259	HEM	C2A-CAA-CBA-CGA
2	C	1259	HEM	C3D-CAD-CBD-CGD
2	A	1258	HEM	C3D-CAD-CBD-CGD
2	B	1259	HEM	CAD-CBD-CGD-O2D
2	C	1259	HEM	CAA-CBA-CGA-O1A
2	A	1258	HEM	CAD-CBD-CGD-O2D
2	B	1259	HEM	CAA-CBA-CGA-O2A
2	A	1258	HEM	CAD-CBD-CGD-O1D

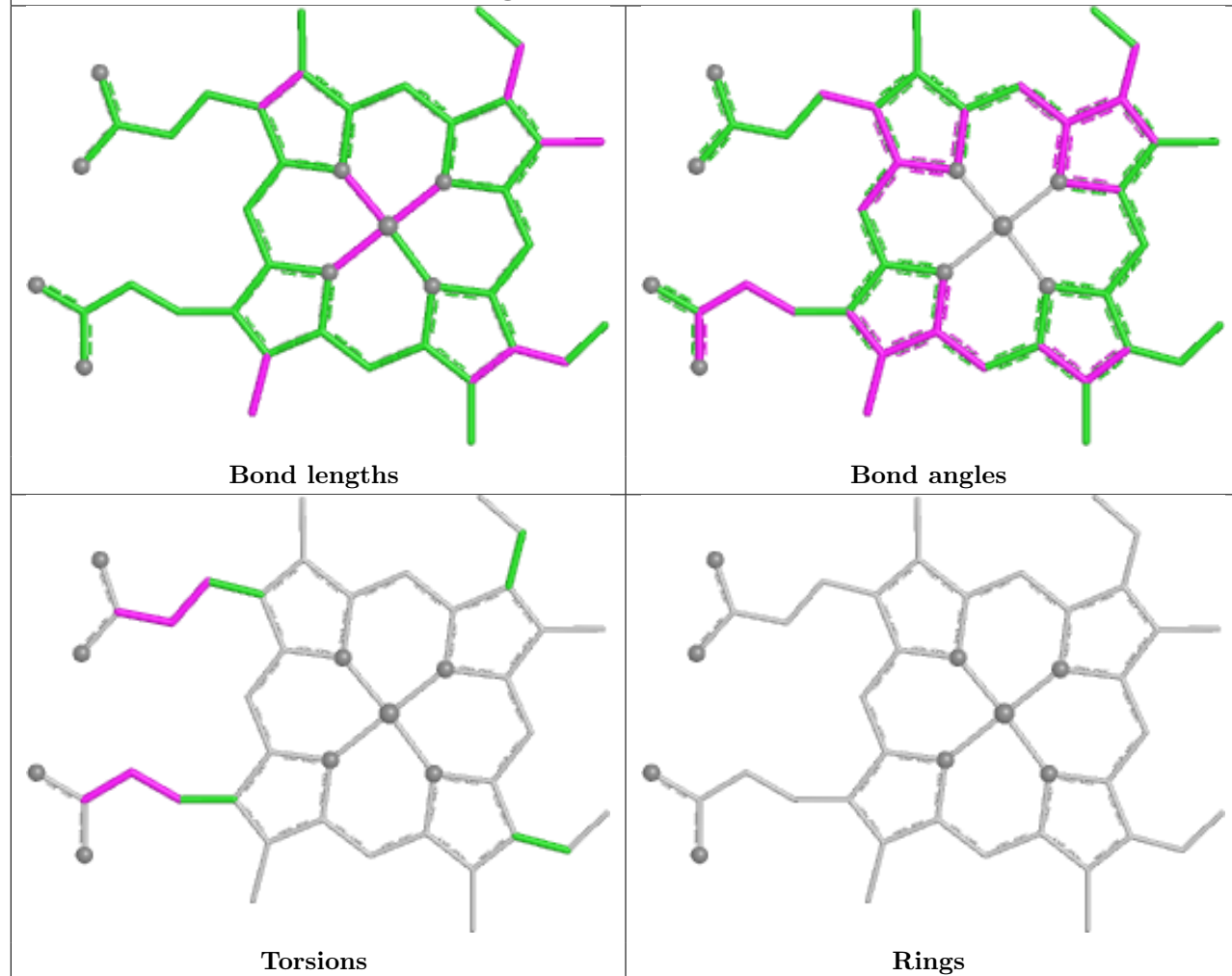
There are no ring outliers.

4 monomers are involved in 12 short contacts:

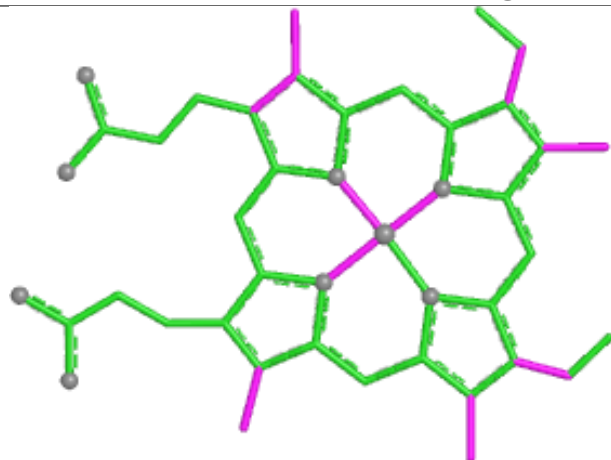
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	1259	HEM	4	0
2	D	1259	HEM	1	0
2	A	1258	HEM	5	0
2	C	1259	HEM	2	0

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.

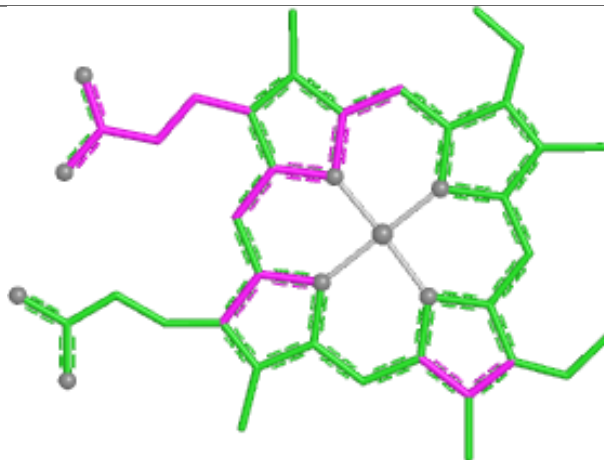
Ligand HEM B 1259



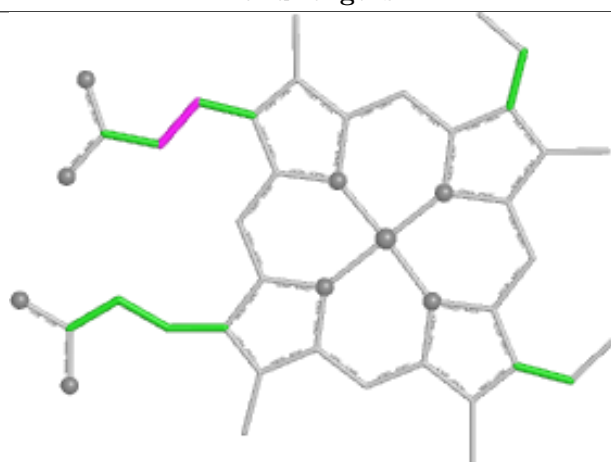
Ligand HEM D 1259



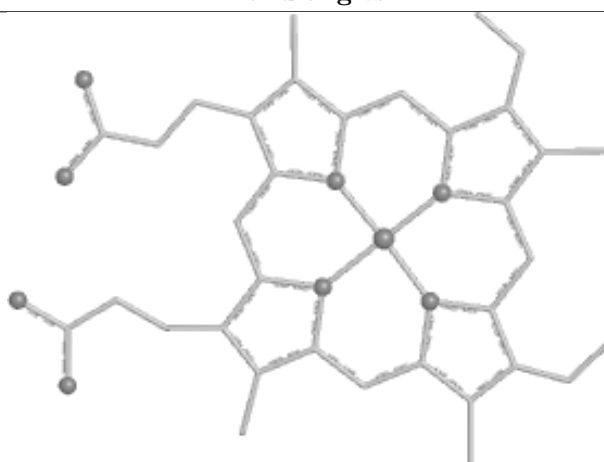
Bond lengths



Bond angles

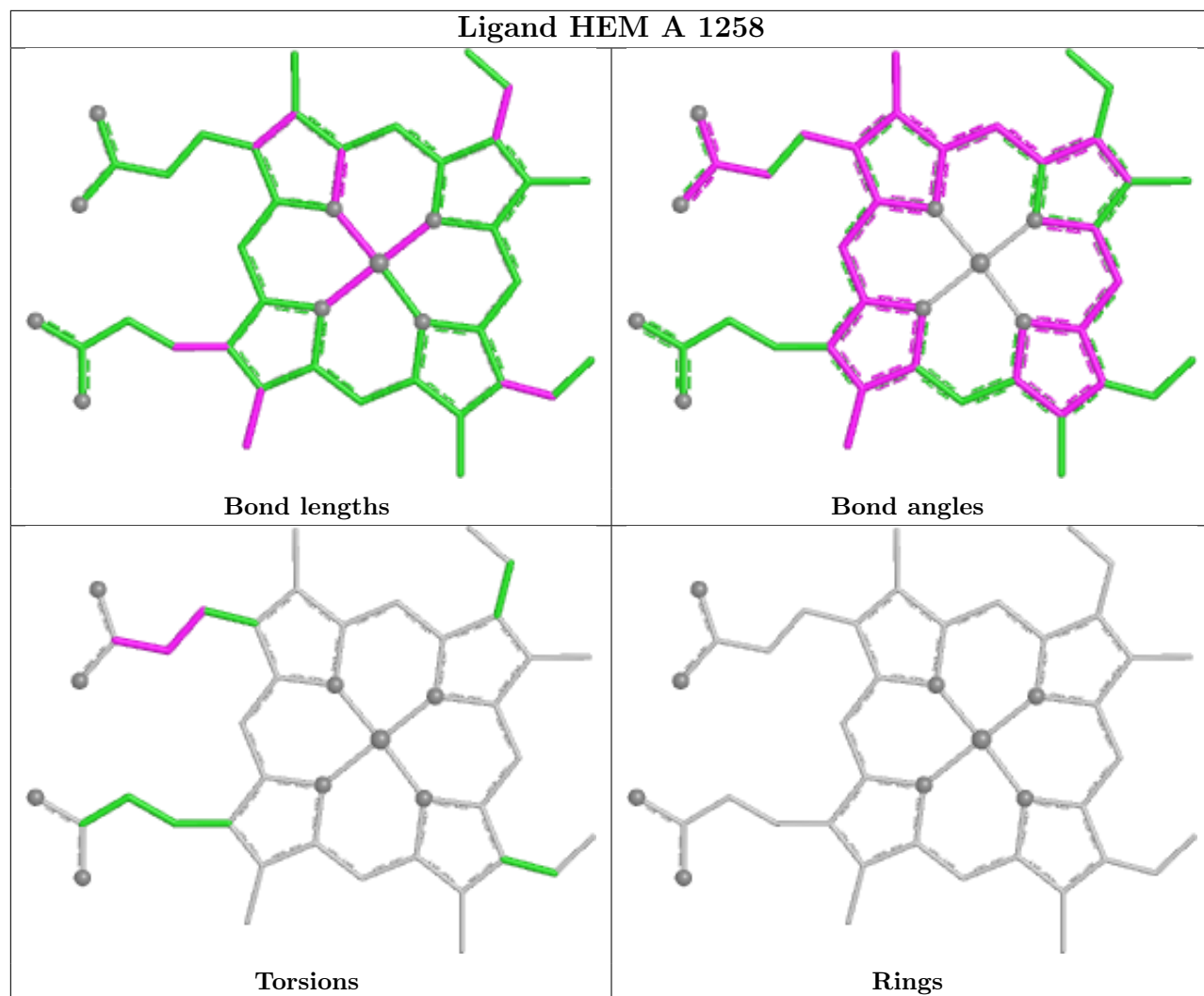


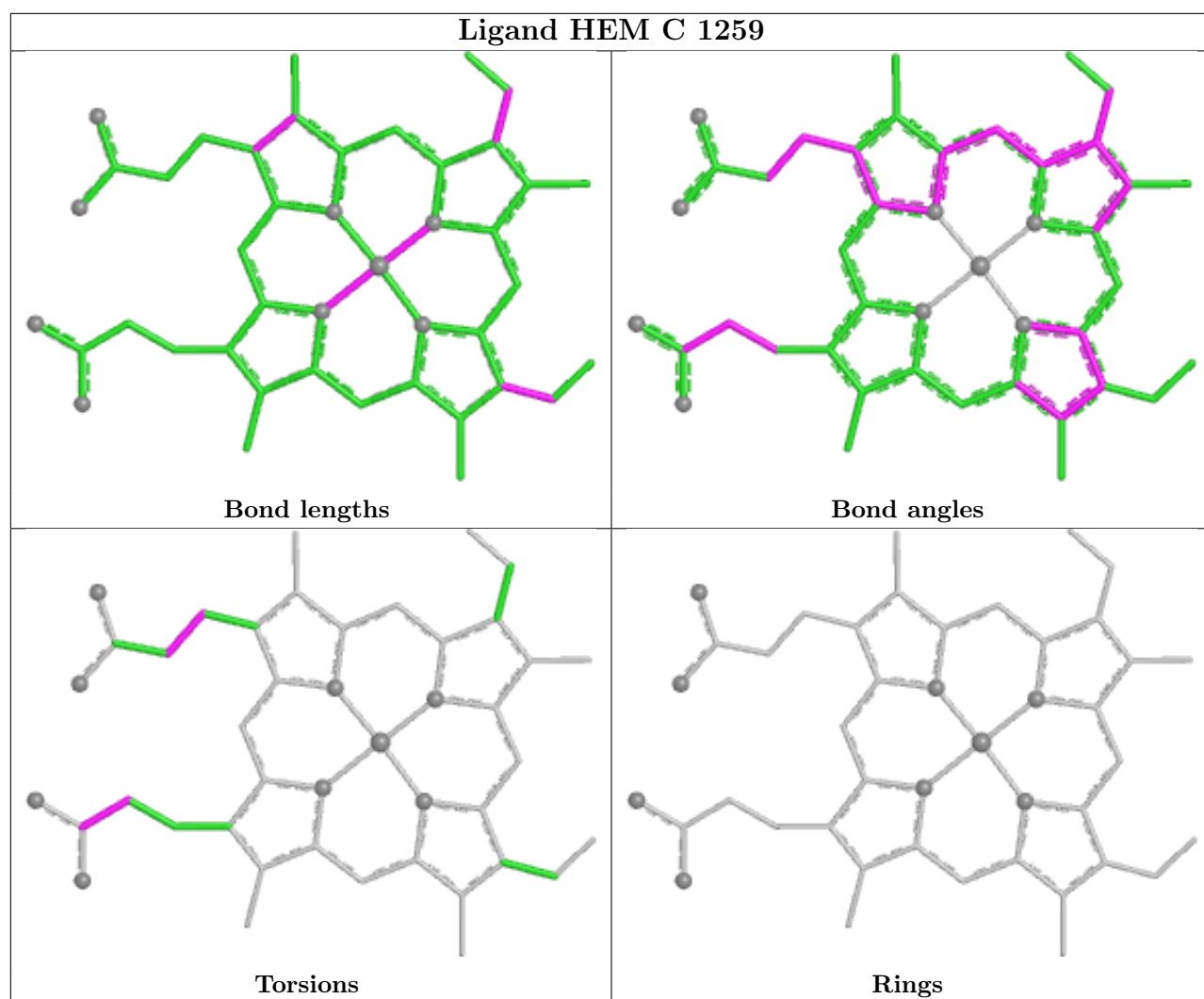
Torsions



Rings

Ligand HEM A 1258





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

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, 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	106/119 (89%)	1.37	23 (21%) 2 2	7, 16, 21, 29	15 (14%)
1	B	107/119 (89%)	1.27	20 (18%) 3 3	9, 15, 20, 31	9 (8%)
1	C	106/119 (89%)	1.33	22 (20%) 2 2	8, 15, 23, 31	6 (5%)
1	D	108/119 (90%)	1.38	27 (25%) 2 1	8, 16, 22, 27	6 (5%)
All	All	427/476 (89%)	1.34	92 (21%) 2 2	7, 16, 22, 31	36 (8%)

All (92) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	258	GLU	9.1
1	B	157	ILE	4.8
1	B	158	VAL	4.3
1	B	230	THR	4.3
1	B	212	ASP	4.2
1	B	152	ILE	4.0
1	D	229	GLY	3.9
1	C	153	PRO	3.8
1	D	212	ASP	3.8
1	A	198	SER	3.7
1	D	210	THR	3.5
1	B	177	GLY	3.5
1	A	157	ILE	3.4
1	D	152	ILE	3.3
1	A	152	ILE	3.3
1	A	245	GLY	3.2
1	A	230	THR	3.2
1	A	168	PHE	3.1
1	C	257	THR	3.1
1	C	256	LEU	3.1
1	D	214	HIS	3.0

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Mol	Chain	Res	Type	RSRZ
1	A	165	ILE	3.0
1	A	231	THR	3.0
1	B	257	THR	3.0
1	D	164	ILE	3.0
1	A	256	LEU	2.9
1	A	158[A]	VAL	2.9
1	C	208	ARG	2.9
1	C	211	SER	2.9
1	A	209	THR	2.9
1	D	151	THR	2.9
1	A	167[A]	LEU	2.8
1	B	155	ALA	2.8
1	A	202	SER	2.8
1	A	210	THR	2.8
1	A	212	ASP	2.7
1	B	159	ILE	2.7
1	D	158	VAL	2.7
1	C	209	THR	2.7
1	C	210	THR	2.7
1	A	211	SER	2.7
1	D	217	GLY	2.7
1	B	245	GLY	2.6
1	C	216	ILE	2.6
1	D	159[A]	ILE	2.6
1	B	258	GLU	2.6
1	B	244	GLY	2.6
1	D	230	THR	2.6
1	D	167	LEU	2.6
1	B	211	SER	2.6
1	D	228	ASP	2.5
1	B	229	GLY	2.5
1	D	256	LEU	2.5
1	C	177	GLY	2.5
1	D	209	THR	2.5
1	C	162	HIS	2.4
1	D	245	GLY	2.4
1	A	199	ARG	2.4
1	B	243	SER	2.3
1	A	195	PRO	2.3
1	D	257	THR	2.3
1	A	184	ILE	2.3
1	C	157	ILE	2.3

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Mol	Chain	Res	Type	RSRZ
1	C	184	ILE	2.3
1	B	252	PHE	2.3
1	D	165	ILE	2.3
1	A	169[A]	SER	2.2
1	B	191	LEU	2.2
1	D	236	LEU	2.2
1	C	232	PHE	2.2
1	A	257	THR	2.2
1	D	172	ALA	2.2
1	D	251	GLY	2.2
1	B	153	PRO	2.2
1	B	165	ILE	2.2
1	C	212	ASP	2.1
1	A	194	GLU	2.1
1	C	254[A]	ARG	2.1
1	A	196	ASP	2.1
1	C	154	ASP	2.1
1	C	206	ARG	2.1
1	B	214	HIS	2.1
1	C	248	TYR	2.1
1	D	157[A]	ILE	2.1
1	D	184	ILE	2.1
1	D	207	TYR	2.1
1	C	167	LEU	2.0
1	D	195	PRO	2.0
1	C	156	MET	2.0
1	C	199	ARG	2.0
1	D	238	ILE	2.0
1	C	183	ALA	2.0

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 oligosaccharides in this entry.

6.4 Ligands

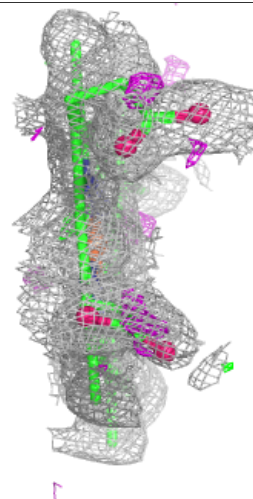
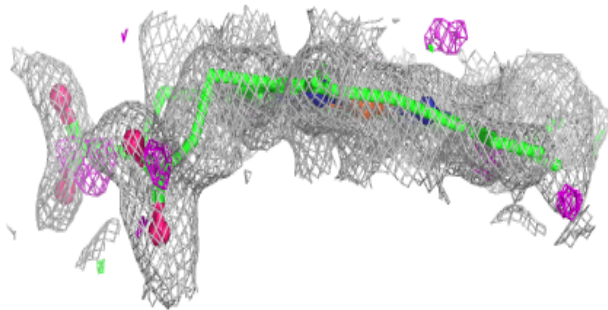
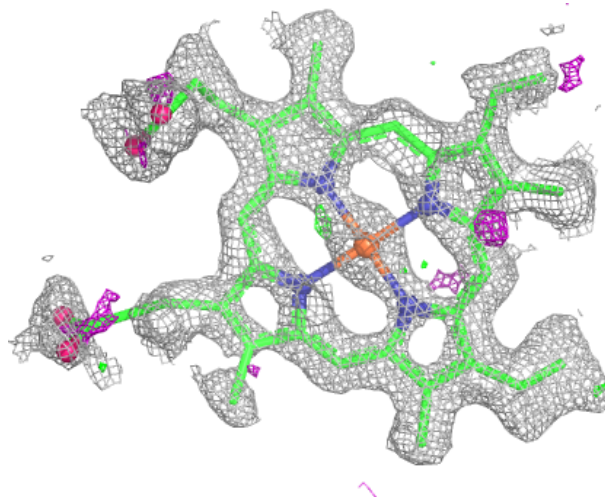
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	CMO	A	1259	2/2	0.89	0.13	6,6,6,12	0
4	NA	C	1261	1/1	0.94	0.07	9,9,9,9	0
2	HEM	A	1258	43/43	0.95	0.10	2,7,20,23	0
2	HEM	C	1259	43/43	0.95	0.09	2,7,15,17	0
2	HEM	D	1259	43/43	0.96	0.09	4,8,18,21	0
3	CMO	C	1260	2/2	0.97	0.04	7,7,7,8	0
3	CMO	D	1260	2/2	0.97	0.12	9,9,9,11	0
2	HEM	B	1259	43/43	0.97	0.07	2,5,16,23	0
3	CMO	B	1260	2/2	0.98	0.07	4,4,4,5	0
5	CL	C	1262	1/1	0.98	0.07	14,14,14,14	0
5	CL	C	1263	1/1	0.99	0.11	16,16,16,16	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.

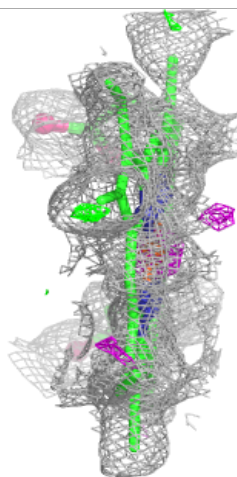
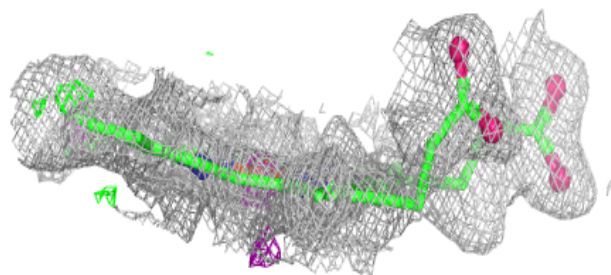
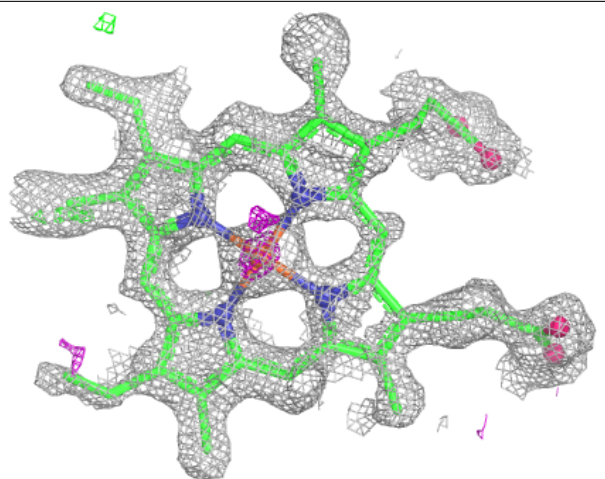
Electron density around HEM A 1258:

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



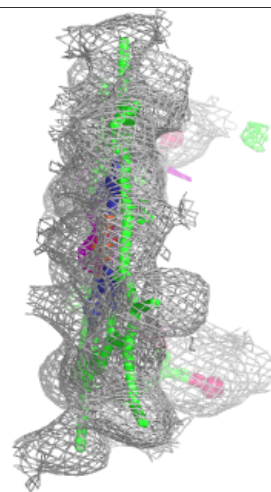
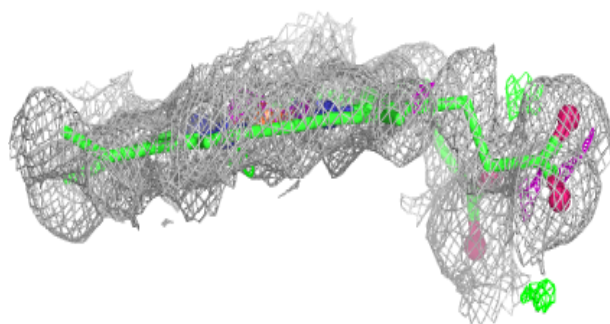
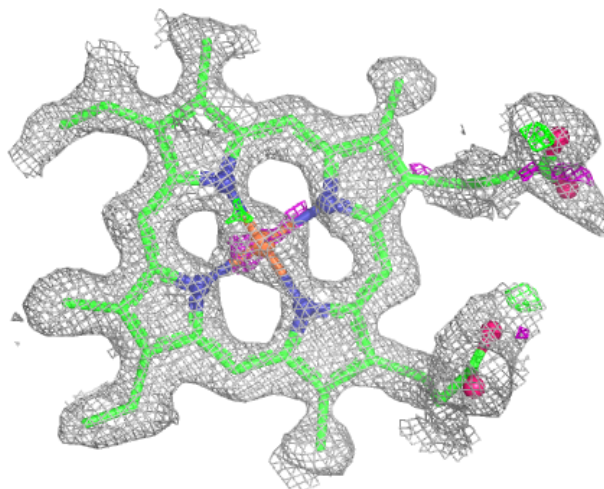
Electron density around HEM C 1259:

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



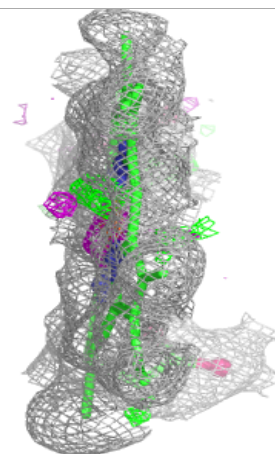
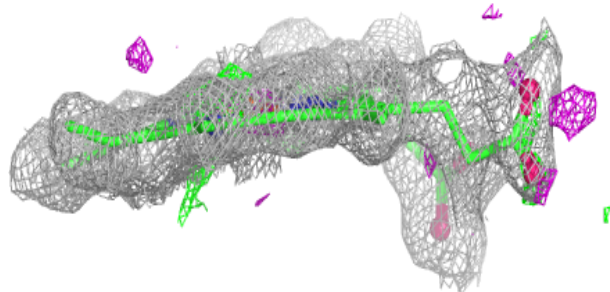
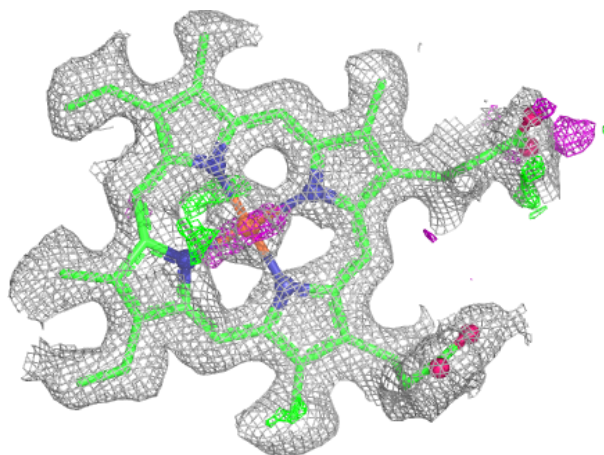
Electron density around HEM D 1259:

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



Electron density around HEM B 1259:

$2mF_o-DF_c$ (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.