



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 06:50 PM UTC

PDB ID : 6AZU / pdb_00006azu
Title : Holo IDO1 crystal structure
Authors : Lewis, H.A.; Yan, C.
Deposited on : 2017-09-13
Resolution : 2.82 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

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

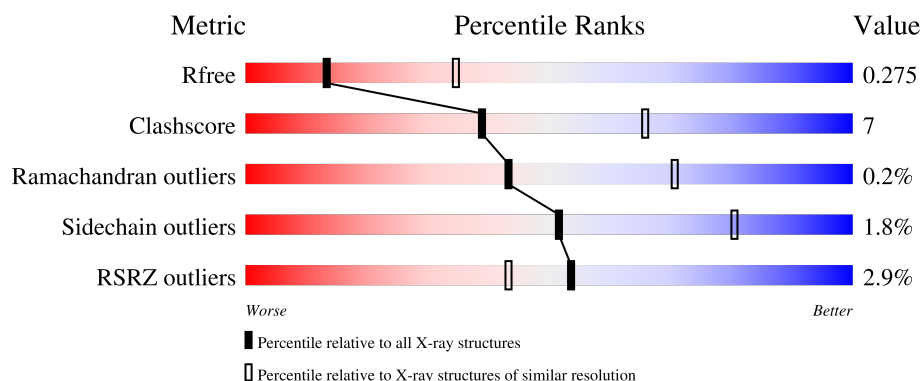
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.82 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	4591 (2.84-2.80)
Clashscore	190562	5010 (2.84-2.80)
Ramachandran outliers	187476	4916 (2.84-2.80)
Sidechain outliers	187428	4918 (2.84-2.80)
RSRZ outliers	180081	4594 (2.84-2.80)

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	402	 2% 76% 15% 8%
1	B	402	 2% 77% 14% 9%
1	C	402	 2% 77% 14% 9%
1	D	402	 4% 75% 16% 8%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 10993 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.

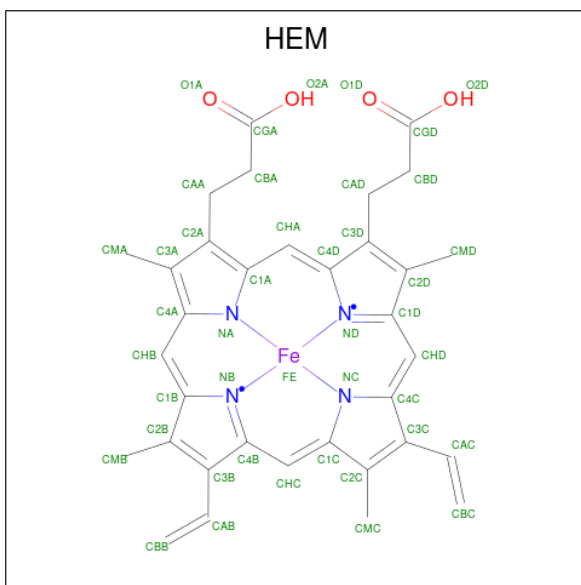
- Molecule 1 is a protein called Indoleamine 2,3-dioxygenase 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	370	Total	C	N	O	S	0	0	0
			2730	1777	458	479	16			
1	B	366	Total	C	N	O	S	0	0	0
			2709	1763	455	475	16			
1	C	366	Total	C	N	O	S	0	0	0
			2686	1746	449	475	16			
1	D	370	Total	C	N	O	S	0	0	0
			2676	1736	450	474	16			

There are 12 discrepancies between the modelled and reference sequences:

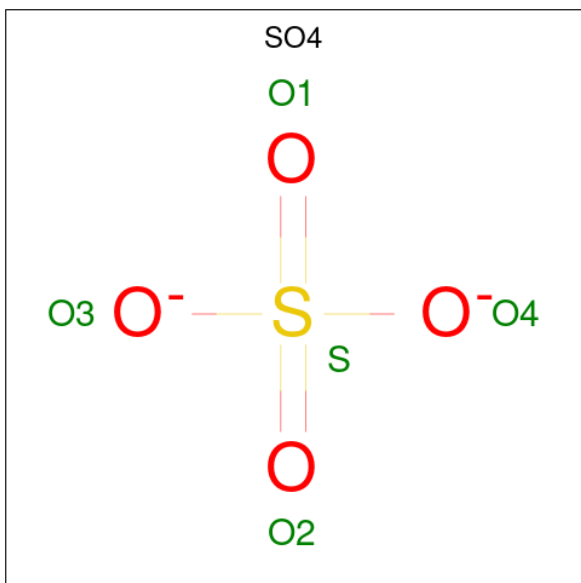
Chain	Residue	Modelled	Actual	Comment	Reference
A	2	GLY	-	expression tag	UNP P14902
A	3	SER	-	expression tag	UNP P14902
A	4	HIS	-	expression tag	UNP P14902
B	2	GLY	-	expression tag	UNP P14902
B	3	SER	-	expression tag	UNP P14902
B	4	HIS	-	expression tag	UNP P14902
C	2	GLY	-	expression tag	UNP P14902
C	3	SER	-	expression tag	UNP P14902
C	4	HIS	-	expression tag	UNP P14902
D	2	GLY	-	expression tag	UNP P14902
D	3	SER	-	expression tag	UNP P14902
D	4	HIS	-	expression tag	UNP P14902

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

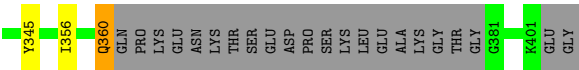


Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total 43	C 34	Fe 1	N 4	O 4	0	0
2	B	1	Total 43	C 34	Fe 1	N 4	O 4	0	0
2	C	1	Total 43	C 34	Fe 1	N 4	O 4	0	0
2	D	1	Total 43	C 34	Fe 1	N 4	O 4	0	0

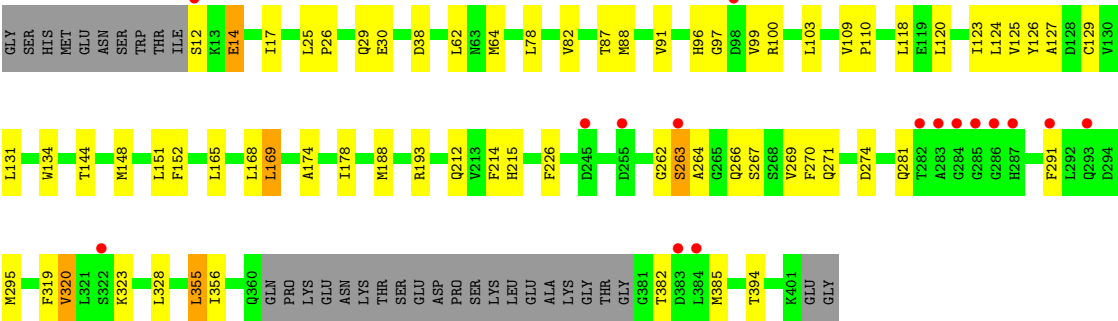
- Molecule 3 is SULFATE ION (CCD ID: SO4) (formula: O_4S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	O	S	0	0
			5	4	1		
3	B	1	Total	O	S	0	0
			5	4	1		
3	C	1	Total	O	S	0	0
			5	4	1		
3	D	1	Total	O	S	0	0
			5	4	1		



● Molecule 1: Indoleamine 2,3-dioxygenase 1



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	119.11Å 177.51Å 101.25Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	34.02 – 2.82 34.02 – 2.82	Depositor EDS
% Data completeness (in resolution range)	99.8 (34.02-2.82) 99.7 (34.02-2.82)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.15	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.05 (at 2.81Å)	Xtriage
Refinement program	PHENIX	Depositor
R, R_{free}	0.211 , 0.272 0.217 , 0.275	Depositor DCC
R_{free} test set	2670 reflections (5.09%)	wwPDB-VP
Wilson B-factor (Å ²)	50.2	Xtriage
Anisotropy	0.872	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 76.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	10993	wwPDB-VP
Average B, all atoms (Å ²)	48.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 14.86% 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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: SO4, 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	0.62	1/2796 (0.0%)	0.97	3/3813 (0.1%)
1	B	0.64	1/2774 (0.0%)	0.94	3/3779 (0.1%)
1	C	0.60	0/2751	0.98	10/3751 (0.3%)
1	D	0.59	0/2742	0.98	8/3742 (0.2%)
All	All	0.61	2/11063 (0.0%)	0.97	24/15085 (0.2%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	109	VAL	CA-CB	5.89	1.57	1.54
1	B	181	ILE	CA-CB	5.56	1.57	1.54

All (24) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	120	LEU	CA-C-N	7.14	124.80	119.66
1	D	120	LEU	C-N-CA	7.14	124.80	119.66
1	C	103	LEU	CA-C-N	6.78	126.74	119.76
1	C	103	LEU	C-N-CA	6.78	126.74	119.76
1	C	120	LEU	CA-C-N	6.63	124.43	119.66
1	C	120	LEU	C-N-CA	6.63	124.43	119.66
1	D	323	LYS	N-CA-C	-6.59	105.86	114.04
1	D	356	ILE	CB-CA-C	-6.33	107.64	114.35
1	A	222	ASN	CA-C-N	6.25	125.74	119.24
1	A	222	ASN	C-N-CA	6.25	125.74	119.24
1	D	355	LEU	N-CA-C	5.92	117.43	110.97
1	C	244	SER	N-CA-C	5.81	118.36	111.33
1	D	382	THR	N-CA-C	-5.78	106.39	113.50
1	B	383	ASP	N-CA-C	-5.64	106.35	113.18
1	A	382	THR	N-CA-C	-5.59	106.62	113.50

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	222	ASN	CA-C-N	5.54	124.73	118.97
1	C	222	ASN	C-N-CA	5.54	124.73	118.97
1	D	96	HIS	CA-C-N	-5.24	117.20	123.44
1	D	96	HIS	C-N-CA	-5.24	117.20	123.44
1	C	27	ASN	CA-C-N	5.17	124.94	119.76
1	C	27	ASN	C-N-CA	5.17	124.94	119.76
1	B	134	TRP	N-CA-C	5.10	116.59	108.79
1	C	13	LYS	N-CA-C	-5.07	105.40	112.45
1	B	95	GLY	N-CA-C	5.01	123.30	115.08

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	2730	0	2512	36	0
1	B	2709	0	2471	34	0
1	C	2686	0	2429	33	0
1	D	2676	0	2377	38	0
2	A	43	0	30	3	0
2	B	43	0	30	1	0
2	C	43	0	30	3	0
2	D	43	0	30	2	0
3	A	5	0	0	0	0
3	B	5	0	0	0	0
3	C	5	0	0	0	0
3	D	5	0	0	0	0
All	All	10993	0	9909	142	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 7.

All (142) 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:C:51:GLU:HG2	1:C:155:ARG:HH21	1.34	0.93
1:C:194:ASP:OD2	1:D:193:ARG:NH2	2.18	0.76
1:B:80:ARG:NH2	1:B:121:PRO:O	2.19	0.76
1:C:176:SER:HA	1:C:179:LYS:HE2	1.66	0.76
1:A:193:ARG:HH12	1:A:197:LEU:HD12	1.55	0.72
1:C:275:VAL:HG13	1:C:311:GLU:HG3	1.74	0.67
1:A:385:MET:HA	1:A:385:MET:HE2	1.75	0.67
1:A:280:GLN:NE2	1:B:293:GLN:OE1	2.28	0.67
2:A:501:HEM:HMB1	2:A:501:HEM:HBB2	1.77	0.66
1:B:275:VAL:HG13	1:B:311:GLU:HG3	1.77	0.66
1:B:385:MET:HA	1:B:385:MET:HE2	1.77	0.65
1:A:275:VAL:HG13	1:A:311:GLU:HG3	1.77	0.64
1:B:88:MET:HE1	1:B:123:ILE:HG13	1.80	0.64
1:D:124:LEU:HD21	1:D:129:CYS:HB3	1.78	0.64
1:D:262:GLY:O	1:D:264:ALA:N	2.28	0.61
1:A:119:GLU:HB2	1:A:301:PRO:HG3	1.82	0.61
1:C:267:SER:OG	1:C:269:VAL:HG12	2.01	0.60
1:B:271:GLN:HA	1:B:274:ASP:HB2	1.85	0.59
1:C:188:MET:HG2	1:C:193:ARG:HH12	1.67	0.59
1:D:88:MET:HE1	1:D:123:ILE:HG13	1.86	0.57
1:C:90:TYR:CE2	1:C:104:PRO:HD3	2.40	0.57
1:D:165:LEU:O	1:D:169:LEU:HD12	2.05	0.57
1:C:127:ALA:HA	1:C:131:LEU:HD12	1.87	0.56
1:C:88:MET:HE1	1:C:123:ILE:HG13	1.88	0.55
1:A:214:PHE:O	1:A:214:PHE:HD2	1.89	0.55
1:B:88:MET:HE3	1:B:124:LEU:H	1.72	0.54
2:B:501:HEM:HMC1	2:B:501:HEM:HBC2	1.89	0.54
1:B:176:SER:HA	1:B:179:LYS:HE2	1.90	0.54
1:D:88:MET:HE3	1:D:124:LEU:H	1.72	0.53
1:C:12:SER:HA	1:C:17:ILE:H	1.73	0.53
2:C:501:HEM:HBC2	2:C:501:HEM:HMC2	1.89	0.53
2:C:501:HEM:HBB2	2:C:501:HEM:HMB1	1.91	0.53
1:C:51:GLU:HG2	1:C:155:ARG:NH2	2.16	0.52
1:A:176:SER:HA	1:A:179:LYS:HE3	1.91	0.52
1:D:134:TRP:CD1	1:D:168:LEU:HD11	2.45	0.52
1:A:25:LEU:HD23	1:A:28:PRO:HB3	1.92	0.51
1:A:204:ALA:O	1:A:208:GLU:HG3	2.10	0.51
1:C:176:SER:HB3	1:C:206:CYS:SG	2.50	0.51
1:A:176:SER:HB3	1:A:206:CYS:SG	2.50	0.51
1:C:356:ILE:HG22	1:C:360:GLN:OE1	2.11	0.51
1:B:350:VAL:HG11	1:B:385:MET:HE3	1.93	0.51
2:A:501:HEM:HBB2	2:A:501:HEM:CMB	2.41	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:80:ARG:HH21	1:C:125:VAL:HG22	1.75	0.50
1:D:30:GLU:HA	1:D:151:LEU:HD21	1.91	0.50
1:D:82:VAL:HG12	1:D:152:PHE:CZ	2.45	0.50
1:A:80:ARG:NH2	1:A:128:ASP:OD2	2.38	0.50
1:D:127:ALA:HA	1:D:131:LEU:HD12	1.93	0.50
1:D:97:GLY:O	1:D:99:VAL:N	2.45	0.49
1:C:218:HIS:NE2	1:C:345:TYR:OH	2.44	0.48
1:B:177:ALA:HB2	1:B:206:CYS:HB2	1.95	0.48
1:B:212:GLN:O	1:B:215:HIS:HB2	2.13	0.48
1:D:29:GLN:O	1:D:78:LEU:HD12	2.14	0.48
1:A:262:GLY:O	1:A:266:GLN:HG3	2.13	0.48
1:B:262:GLY:O	1:B:266:GLN:HG3	2.13	0.48
1:A:357:PRO:C	1:A:359:SER:H	2.22	0.48
1:C:326:ALA:O	1:C:330:GLU:HB2	2.14	0.48
1:D:271:GLN:HA	1:D:274:ASP:HB2	1.95	0.48
1:A:271:GLN:HA	1:A:274:ASP:HB2	1.96	0.47
1:D:38:ASP:HB3	1:D:62:LEU:HD11	1.95	0.47
1:D:174:ALA:O	1:D:178:ILE:HG13	2.15	0.47
1:D:109:VAL:HB	1:D:110:PRO:HD3	1.95	0.47
1:A:311:GLU:HG2	1:B:293:GLN:NE2	2.30	0.47
1:B:222:ASN:OD1	1:B:225:ALA:N	2.36	0.47
1:A:267:SER:OG	1:A:269:VAL:HG12	2.15	0.47
1:B:119:GLU:HB2	1:B:301:PRO:HG3	1.96	0.47
1:C:148:MET:HE1	1:C:168:LEU:HD12	1.96	0.47
1:B:237:TRP:HB3	1:B:243:LEU:HD22	1.97	0.46
2:D:501:HEM:HMC1	2:D:501:HEM:HBC2	1.97	0.46
1:A:125:VAL:HG12	1:A:127:ALA:H	1.80	0.46
1:C:218:HIS:HE2	1:C:345:TYR:HH	1.64	0.46
1:D:212:GLN:O	1:D:215:HIS:HB2	2.15	0.46
1:D:134:TRP:CG	1:D:168:LEU:HD11	2.51	0.46
1:B:352:LYS:HD3	1:B:353:TYR:CE1	2.51	0.46
1:D:25:LEU:HA	1:D:26:PRO:HD3	1.78	0.46
1:D:126:TYR:CB	1:D:266:GLN:HB2	2.46	0.45
1:A:356:ILE:HB	1:A:357:PRO:HD3	1.98	0.45
1:A:312:SER:O	1:B:297:ARG:HG2	2.16	0.45
1:B:349:ILE:HD13	1:B:349:ILE:HA	1.85	0.45
1:B:124:LEU:HD21	1:B:129:CYS:HB3	1.99	0.45
1:A:115:SER:HB3	1:A:120:LEU:O	2.17	0.44
1:D:12:SER:C	1:D:14:GLU:H	2.24	0.44
1:D:100:ARG:HA	1:D:100:ARG:HD3	1.73	0.44
1:A:213:VAL:O	1:A:216:GLN:HB2	2.16	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:118:LEU:HD23	1:D:118:LEU:HA	1.72	0.44
1:D:263:SER:HA	1:D:266:GLN:OE1	2.18	0.44
1:D:267:SER:OG	1:D:269:VAL:HG12	2.17	0.44
1:C:185:PHE:HA	1:C:188:MET:HE2	1.99	0.44
1:D:144:THR:O	1:D:148:MET:HG3	2.17	0.44
1:A:269:VAL:HG13	1:A:270:PHE:CD1	2.53	0.44
1:C:82:VAL:HG12	1:C:152:PHE:CZ	2.53	0.44
1:A:24:ALA:HA	1:A:131:LEU:HD22	2.00	0.44
1:B:82:VAL:HG12	1:B:152:PHE:CZ	2.52	0.44
1:B:78:LEU:O	1:B:82:VAL:HG13	2.17	0.43
1:C:296:ARG:HA	1:C:299:MET:SD	2.58	0.43
1:C:356:ILE:O	1:C:360:GLN:HG2	2.18	0.43
1:D:355:LEU:HD11	1:D:385:MET:CG	2.47	0.43
1:B:24:ALA:HA	1:B:131:LEU:HD22	2.00	0.43
1:C:174:ALA:O	1:C:178:ILE:HG13	2.18	0.43
1:B:80:ARG:NH1	1:B:128:ASP:OD2	2.51	0.43
1:C:270:PHE:O	1:C:274:ASP:HB2	2.18	0.43
1:D:226:PHE:CZ	2:D:501:HEM:HMB3	2.53	0.43
1:D:320:VAL:HG22	1:D:328:LEU:HB3	1.99	0.43
1:B:127:ALA:HA	1:B:131:LEU:HD12	1.99	0.43
1:B:193:ARG:HD3	1:B:193:ARG:C	2.44	0.43
1:A:222:ASN:HA	1:A:223:PRO:HD3	1.88	0.43
1:A:311:GLU:HG2	1:B:293:GLN:HE21	1.84	0.43
1:C:88:MET:HE1	1:C:237:TRP:CH2	2.54	0.43
1:C:201:LEU:HD23	1:C:201:LEU:HA	1.74	0.43
1:D:281:GLN:HA	1:D:394:THR:HG21	1.99	0.43
1:C:317:ARG:HA	1:C:320:VAL:HG22	2.00	0.42
1:A:148:MET:HE1	1:A:168:LEU:HD23	2.01	0.42
1:C:214:PHE:HZ	2:C:501:HEM:HHC	1.84	0.42
1:A:78:LEU:O	1:A:82:VAL:HG13	2.19	0.42
1:D:188:MET:HE3	1:D:319:PHE:HB2	2.02	0.42
1:B:121:PRO:HG2	1:B:298:TYR:CD2	2.54	0.42
1:B:23:PHE:O	1:B:131:LEU:HD13	2.20	0.42
1:A:88:MET:HE3	1:A:124:LEU:H	1.84	0.42
1:C:180:VAL:HG12	1:C:180:VAL:O	2.20	0.42
1:D:126:TYR:HB2	1:D:266:GLN:HB2	2.02	0.42
1:A:103:LEU:HA	1:A:104:PRO:HD3	1.96	0.41
1:D:269:VAL:HG13	1:D:270:PHE:CD1	2.55	0.41
1:D:291:PHE:O	1:D:295:MET:HG2	2.19	0.41
1:D:12:SER:C	1:D:14:GLU:N	2.79	0.41
1:A:237:TRP:O	1:A:258:GLU:HG2	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:80:ARG:HD3	1:B:128:ASP:OD1	2.20	0.41
1:A:265:GLY:HA2	2:A:501:HEM:C4D	2.56	0.41
1:B:90:TYR:CE2	1:B:104:PRO:HD3	2.55	0.41
1:A:223:PRO:HG3	1:A:352:LYS:HE2	2.02	0.41
1:C:12:SER:HB2	1:C:17:ILE:O	2.20	0.41
1:D:87:THR:HG23	1:D:103:LEU:HD13	2.02	0.41
1:C:193:ARG:HD2	1:C:319:PHE:CZ	2.56	0.41
1:C:247:LEU:HD23	1:C:249:TYR:CZ	2.55	0.41
1:A:347:LEU:HD21	1:A:388:LEU:HB2	2.03	0.41
1:C:145:TYR:OH	1:C:161:LYS:HG2	2.21	0.41
1:D:214:PHE:O	1:D:214:PHE:HD2	2.03	0.41
1:B:56:ARG:HD3	1:B:100:ARG:CZ	2.50	0.41
1:B:188:MET:HE3	1:B:319:PHE:HB2	2.03	0.41
1:A:192:GLU:HB3	1:A:195:THR:OG1	2.21	0.40
1:A:165:LEU:HA	1:A:165:LEU:HD23	1.77	0.40
1:A:310:LEU:HD23	1:A:310:LEU:HA	1.82	0.40
1:B:181:ILE:HD13	1:B:276:LEU:HD22	2.04	0.40
1:D:91:VAL:O	1:D:99:VAL:HA	2.21	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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	366/402 (91%)	350 (96%)	14 (4%)	2 (0%)	24	53
1	B	360/402 (90%)	349 (97%)	11 (3%)	0	100	100
1	C	360/402 (90%)	347 (96%)	13 (4%)	0	100	100
1	D	366/402 (91%)	352 (96%)	13 (4%)	1 (0%)	36	64
All	All	1452/1608 (90%)	1398 (96%)	51 (4%)	3 (0%)	43	71

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	263	SER
1	A	13	LYS
1	A	285	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	245/348 (70%)	241 (98%)	4 (2%)	55	82
1	B	240/348 (69%)	234 (98%)	6 (2%)	42	74
1	C	237/348 (68%)	236 (100%)	1 (0%)	84	94
1	D	227/348 (65%)	221 (97%)	6 (3%)	40	73
All	All	949/1392 (68%)	932 (98%)	17 (2%)	51	80

All (17) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	125	VAL
1	A	191	GLN
1	A	320	VAL
1	A	356	ILE
1	B	64	MET
1	B	82	VAL
1	B	125	VAL
1	B	268	SER
1	B	320	VAL
1	B	323	LYS
1	C	360	GLN
1	D	14	GLU
1	D	17	ILE
1	D	64	MET
1	D	125	VAL
1	D	169	LEU
1	D	320	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (12)

such sidechains are listed below:

Mol	Chain	Res	Type
1	A	73	HIS
1	A	220	HIS
1	A	242	GLN
1	A	293	GLN
1	A	313	ASN
1	A	348	GLN
1	B	73	HIS
1	B	215	HIS
1	C	212	GLN
1	C	216	GLN
1	C	305	ASN
1	D	360	GLN

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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

8 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	SO4	D	502	-	4,4,4	0.21	0	6,6,6	0.19	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	SO4	A	502	-	4,4,4	0.21	0	6,6,6	0.19	0
2	HEM	C	501	1	50,50,50	1.69	9 (18%)	67,82,82	1.10	4 (5%)
2	HEM	B	501	1	50,50,50	1.69	5 (10%)	67,82,82	1.20	10 (14%)
2	HEM	A	501	1	50,50,50	1.74	11 (22%)	67,82,82	1.77	15 (22%)
2	HEM	D	501	1	50,50,50	1.79	10 (20%)	67,82,82	1.12	3 (4%)
3	SO4	C	502	-	4,4,4	0.26	0	6,6,6	0.13	0
3	SO4	B	502	-	4,4,4	0.24	0	6,6,6	0.11	0

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	C	501	1	-	7/14/54/54	-
2	HEM	A	501	1	-	6/14/54/54	-
2	HEM	D	501	1	-	4/14/54/54	-
2	HEM	B	501	1	-	3/14/54/54	-

All (35) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	501	HEM	C3D-C2D	7.62	1.53	1.36
2	C	501	HEM	C3D-C2D	7.46	1.52	1.36
2	B	501	HEM	C3D-C2D	7.07	1.52	1.36
2	A	501	HEM	C3D-C2D	6.51	1.50	1.36
2	B	501	HEM	FE-NA	4.21	2.09	1.95
2	A	501	HEM	FE-NA	4.14	2.08	1.95
2	D	501	HEM	FE-NB	4.10	2.07	1.94
2	A	501	HEM	FE-NB	-3.61	1.83	1.94
2	D	501	HEM	FE-ND	3.42	2.05	1.94
2	D	501	HEM	CAB-C3B	3.20	1.55	1.47
2	D	501	HEM	FE-NA	3.19	2.05	1.95
2	B	501	HEM	CAC-C3C	3.17	1.55	1.47
2	C	501	HEM	CAB-C3B	3.14	1.55	1.47
2	B	501	HEM	CAB-C3B	3.08	1.55	1.47
2	C	501	HEM	CAC-C3C	3.04	1.55	1.47
2	C	501	HEM	CMC-C2C	2.88	1.56	1.50
2	A	501	HEM	FE-NC	2.62	2.03	1.95
2	A	501	HEM	CAC-C3C	2.61	1.54	1.47

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	501	HEM	C1D-ND	2.60	1.44	1.38
2	D	501	HEM	CAC-C3C	2.55	1.54	1.47
2	A	501	HEM	CAB-C3B	2.54	1.54	1.47
2	C	501	HEM	FE-ND	2.51	2.02	1.94
2	C	501	HEM	FE-NC	2.45	2.03	1.95
2	A	501	HEM	CMC-C2C	2.42	1.55	1.50
2	D	501	HEM	CMA-C3A	2.38	1.55	1.50
2	D	501	HEM	CMC-C2C	2.28	1.55	1.50
2	D	501	HEM	CMB-C2B	2.16	1.55	1.50
2	B	501	HEM	CMC-C2C	2.16	1.55	1.50
2	C	501	HEM	CMA-C3A	2.13	1.55	1.50
2	A	501	HEM	CMA-C3A	2.11	1.55	1.50
2	C	501	HEM	CMB-C2B	2.11	1.55	1.50
2	A	501	HEM	CMD-C2D	2.08	1.55	1.50
2	A	501	HEM	FE-ND	-2.05	1.88	1.94
2	C	501	HEM	FE-NB	2.03	2.01	1.94
2	D	501	HEM	CMD-C2D	2.02	1.54	1.50

All (32) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	501	HEM	CHD-C1D-ND	5.63	130.48	124.42
2	A	501	HEM	CHB-C1B-NB	4.77	130.27	124.37
2	A	501	HEM	CHD-C1D-C2D	-4.69	117.62	125.03
2	D	501	HEM	C4D-ND-C1D	4.45	110.48	105.21
2	C	501	HEM	C4D-ND-C1D	4.02	109.97	105.21
2	A	501	HEM	CHC-C4B-NB	3.50	128.19	124.42
2	A	501	HEM	CAD-C3D-C4D	2.87	129.71	124.70
2	B	501	HEM	CHD-C1D-ND	2.81	127.44	124.42
2	B	501	HEM	C4D-ND-C1D	2.76	108.48	105.21
2	B	501	HEM	CHB-C1B-NB	2.74	127.76	124.37
2	A	501	HEM	C3D-C4D-ND	2.56	112.98	110.17
2	D	501	HEM	CHD-C4C-NC	2.54	127.22	124.45
2	A	501	HEM	CHA-C4D-C3D	-2.45	120.71	125.23
2	B	501	HEM	O2D-CGD-CBD	2.45	121.73	114.00
2	A	501	HEM	C3B-C2B-C1B	2.43	108.24	106.41
2	C	501	HEM	CBA-CAA-C2A	2.42	119.24	112.53
2	C	501	HEM	C3B-C2B-C1B	2.41	108.22	106.41
2	A	501	HEM	CHB-C1B-C2B	-2.36	120.25	126.95
2	A	501	HEM	O1D-CGD-CBD	-2.34	115.68	123.09
2	A	501	HEM	CHC-C1C-NC	-2.24	122.02	124.45
2	D	501	HEM	C3B-C2B-C1B	2.22	108.08	106.41

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	501	HEM	O2D-CGD-CBD	2.18	120.90	114.00
2	A	501	HEM	CAA-C2A-C3A	2.18	131.97	127.07
2	B	501	HEM	CMD-C2D-C1D	2.16	128.41	125.03
2	A	501	HEM	CBA-CAA-C2A	2.15	118.49	112.53
2	B	501	HEM	CHA-C4D-ND	2.12	126.99	124.37
2	B	501	HEM	C1C-CHC-C4B	-2.11	121.53	126.02
2	A	501	HEM	C1C-CHC-C4B	-2.10	121.55	126.02
2	B	501	HEM	CAD-C3D-C4D	2.09	128.34	124.70
2	B	501	HEM	CBD-CAD-C3D	-2.06	106.83	112.53
2	C	501	HEM	CMD-C2D-C1D	2.06	128.25	125.03
2	B	501	HEM	C3B-C2B-C1B	2.04	107.94	106.41

There are no chirality outliers.

All (20) torsion outliers are listed below:

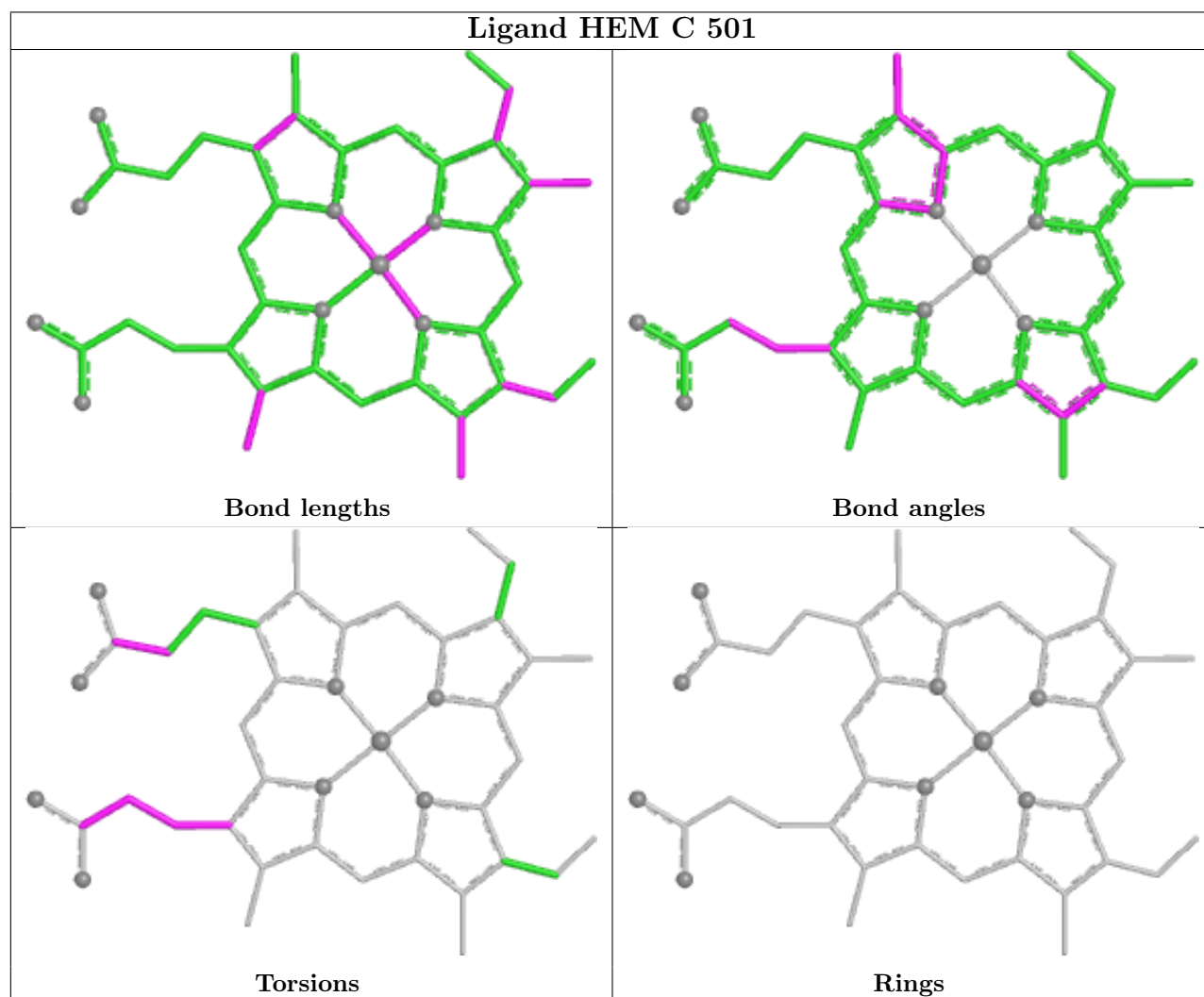
Mol	Chain	Res	Type	Atoms
2	A	501	HEM	C1A-C2A-CAA-CBA
2	A	501	HEM	C3A-C2A-CAA-CBA
2	C	501	HEM	C1A-C2A-CAA-CBA
2	D	501	HEM	C1A-C2A-CAA-CBA
2	D	501	HEM	C3A-C2A-CAA-CBA
2	C	501	HEM	C2A-CAA-CBA-CGA
2	C	501	HEM	C3A-C2A-CAA-CBA
2	A	501	HEM	C2A-CAA-CBA-CGA
2	A	501	HEM	C3D-CAD-CBD-CGD
2	C	501	HEM	CAD-CBD-CGD-O1D
2	A	501	HEM	CAA-CBA-CGA-O2A
2	C	501	HEM	CAD-CBD-CGD-O2D
2	A	501	HEM	CAA-CBA-CGA-O1A
2	B	501	HEM	C3D-CAD-CBD-CGD
2	C	501	HEM	CAA-CBA-CGA-O2A
2	C	501	HEM	CAA-CBA-CGA-O1A
2	D	501	HEM	C3D-CAD-CBD-CGD
2	D	501	HEM	CAD-CBD-CGD-O1D
2	B	501	HEM	CAD-CBD-CGD-O2D
2	B	501	HEM	CAD-CBD-CGD-O1D

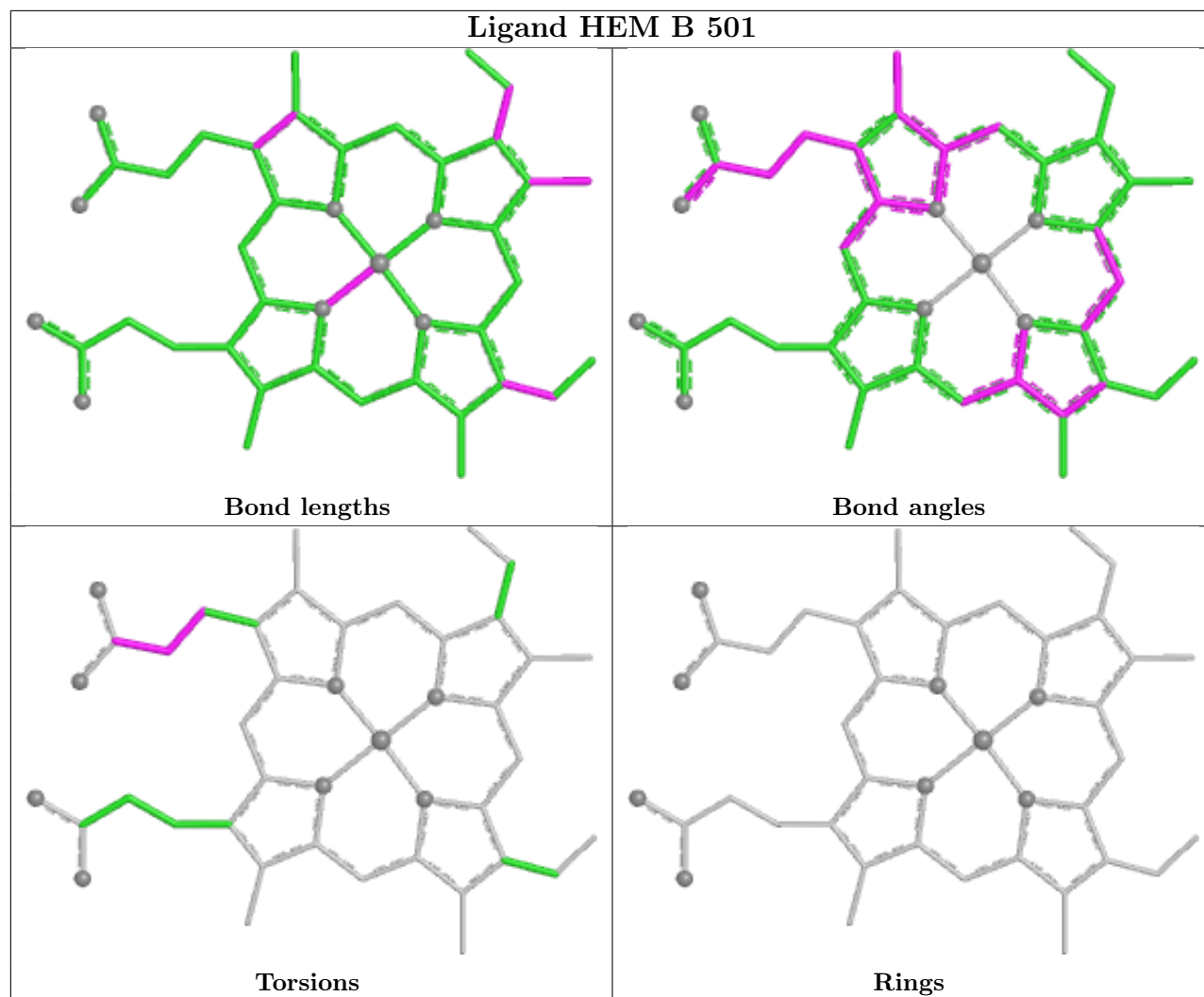
There are no ring outliers.

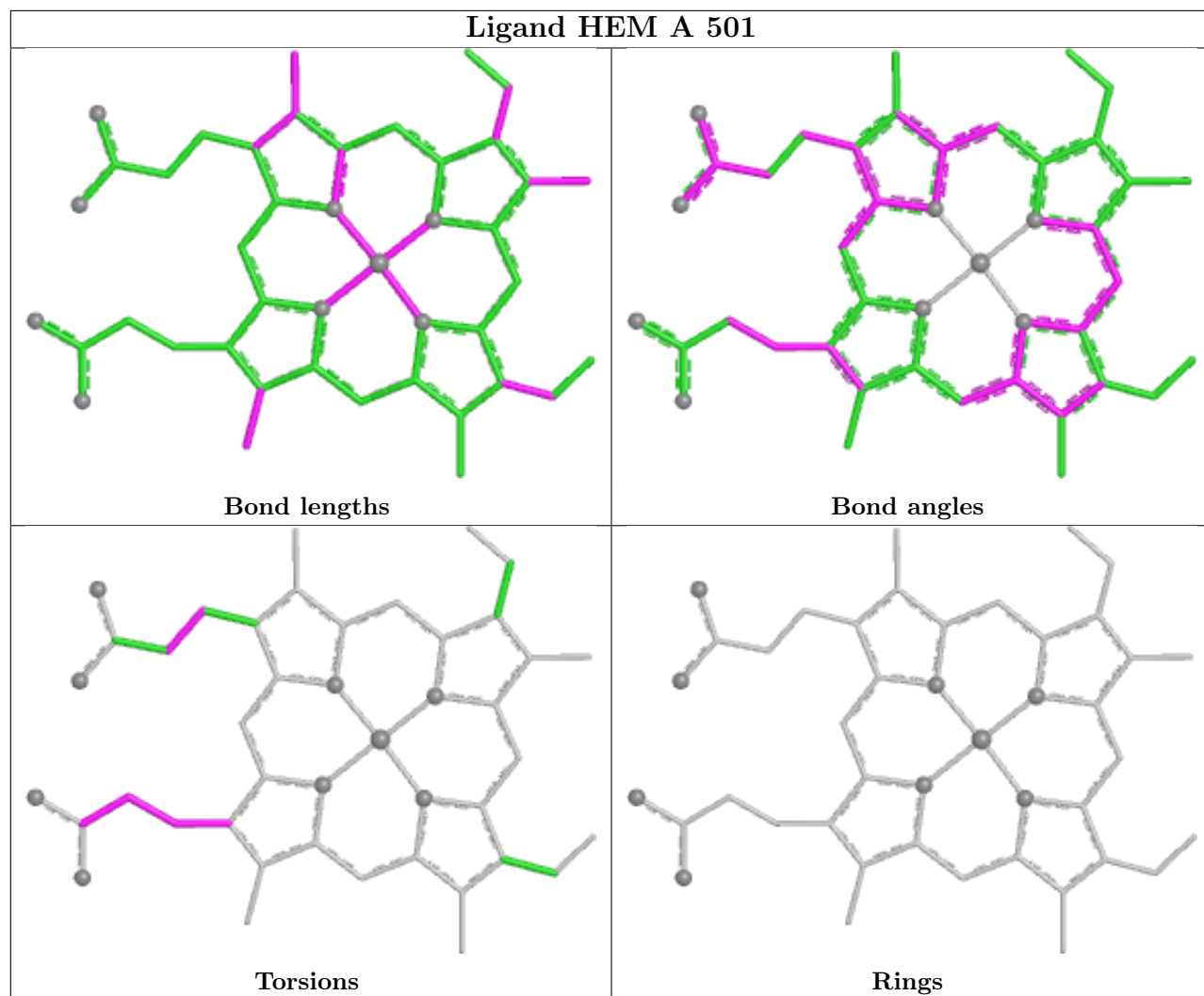
4 monomers are involved in 9 short contacts:

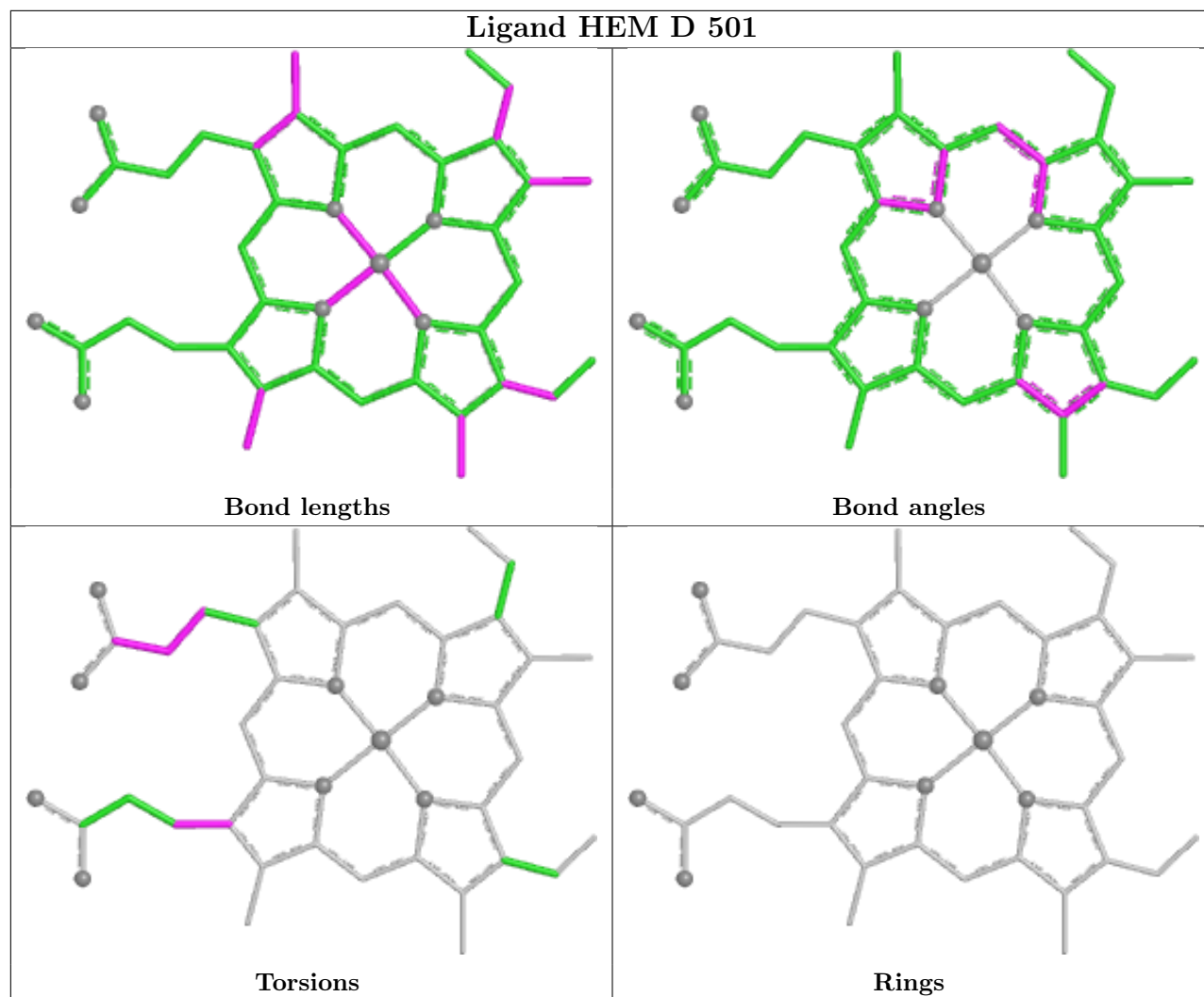
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	501	HEM	3	0
2	B	501	HEM	1	0
2	A	501	HEM	3	0
2	D	501	HEM	2	0

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.









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	370/402 (92%)	-0.04	10 (2%) 56 46	28, 45, 74, 161	0
1	B	366/402 (91%)	-0.04	8 (2%) 62 52	28, 43, 68, 107	0
1	C	366/402 (91%)	0.07	8 (2%) 62 52	29, 47, 75, 122	0
1	D	370/402 (92%)	0.13	16 (4%) 40 31	31, 51, 77, 129	0
All	All	1472/1608 (91%)	0.03	42 (2%) 53 43	28, 47, 74, 161	0

All (42) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	382	THR	4.7
1	C	193	ARG	4.7
1	A	284	GLY	4.4
1	B	287	HIS	4.3
1	A	286	GLY	4.3
1	A	285	GLY	4.2
1	C	319	PHE	4.1
1	B	12	SER	4.0
1	D	285	GLY	3.9
1	D	282	THR	3.8
1	A	12	SER	3.4
1	B	98	ASP	3.3
1	C	324	GLY	3.2
1	A	360	GLN	3.0
1	A	283	ALA	2.9
1	C	287	HIS	2.9
1	D	287	HIS	2.9
1	D	293	GLN	2.9
1	D	12	SER	2.9
1	A	312	SER	2.8
1	D	384	LEU	2.8

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Mol	Chain	Res	Type	RSRZ
1	D	245	ASP	2.7
1	D	98	ASP	2.7
1	D	263	SER	2.5
1	C	291	PHE	2.4
1	D	283	ALA	2.4
1	B	19	GLU	2.3
1	C	214	PHE	2.3
1	D	286	GLY	2.3
1	B	31	ASN	2.2
1	D	255	ASP	2.2
1	C	245	ASP	2.2
1	B	95	GLY	2.2
1	A	99	VAL	2.1
1	B	323	LYS	2.1
1	D	291	PHE	2.1
1	D	322	SER	2.1
1	A	27	ASN	2.1
1	D	284	GLY	2.1
1	A	382	THR	2.1
1	D	383	ASP	2.1
1	C	194	ASP	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(Å ²)	Q<0.9
3	SO4	A	502	5/5	0.72	0.18	117,120,123,127	0
3	SO4	D	502	5/5	0.77	0.14	102,106,110,114	0

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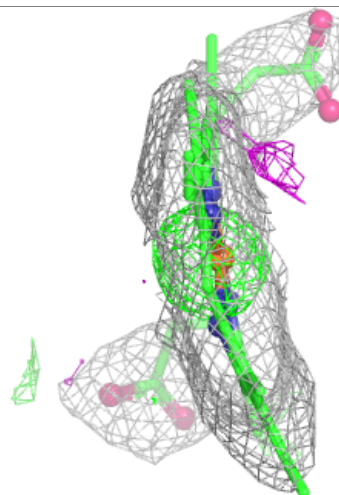
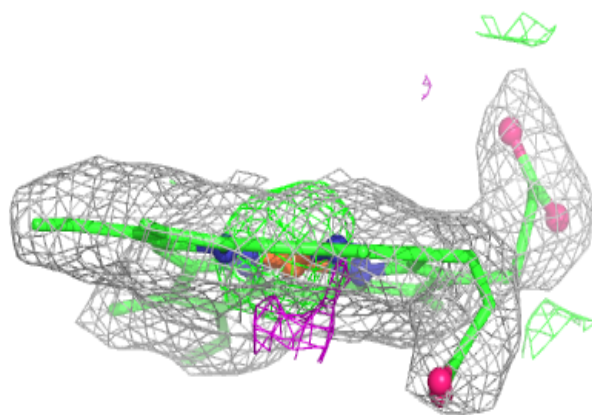
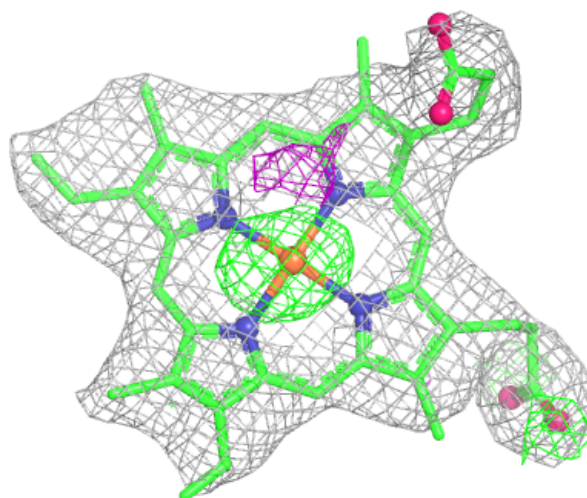
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	SO4	B	502	5/5	0.78	0.15	132,132,136,137	0
3	SO4	C	502	5/5	0.88	0.12	100,100,103,103	0
2	HEM	C	501	43/43	0.91	0.12	22,45,69,108	0
2	HEM	A	501	43/43	0.93	0.11	3,28,75,83	0
2	HEM	D	501	43/43	0.95	0.10	13,49,79,86	0
2	HEM	B	501	43/43	0.97	0.09	9,28,49,53	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.

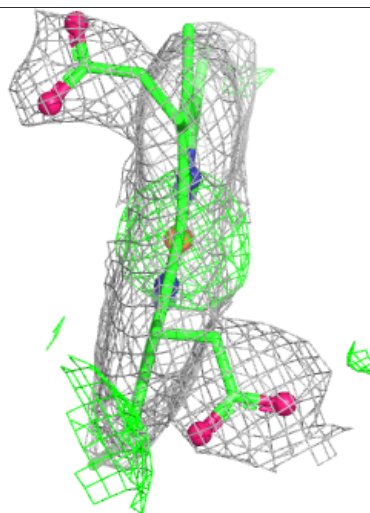
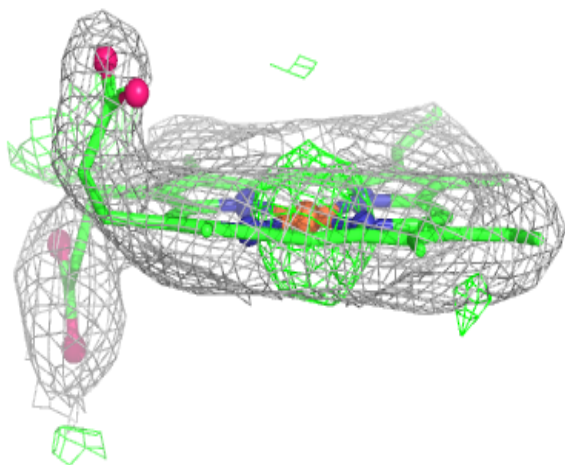
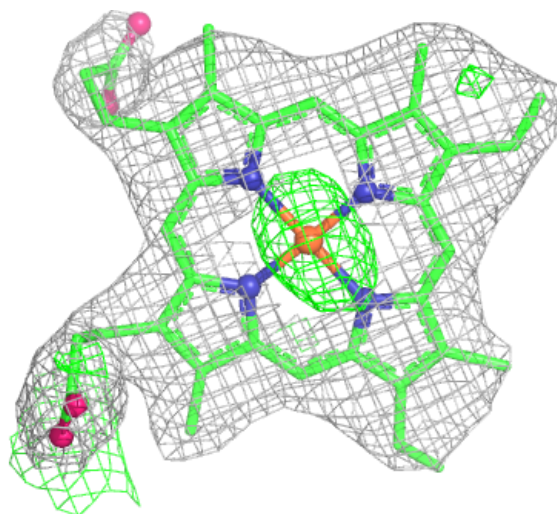
Electron density around HEM C 501:

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



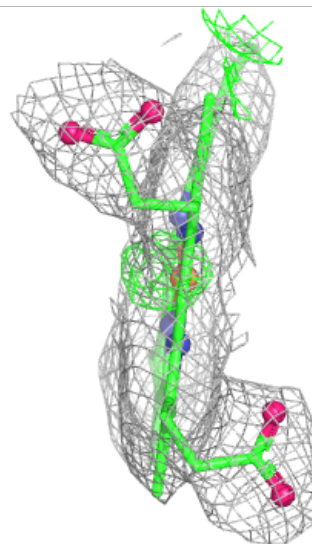
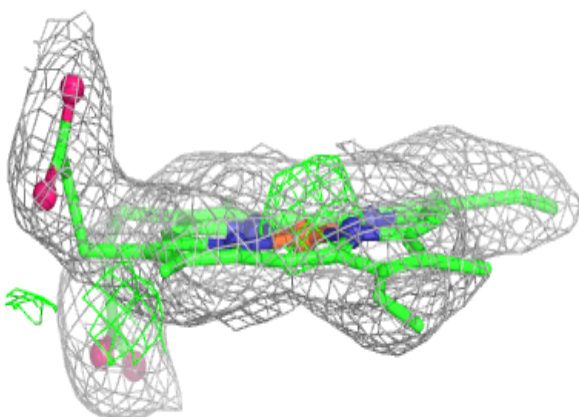
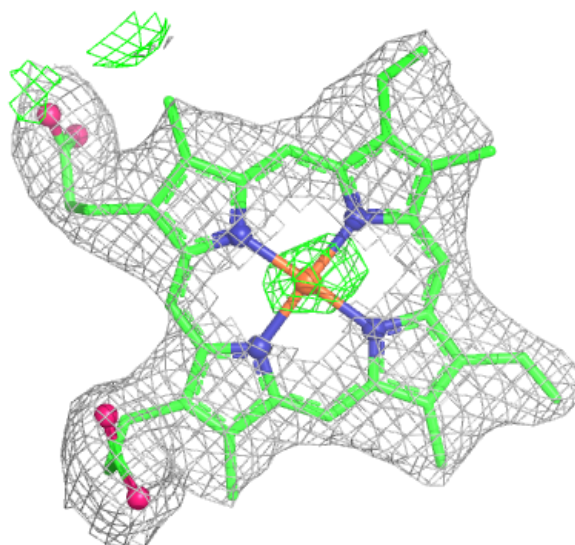
Electron density around HEM A 501:

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



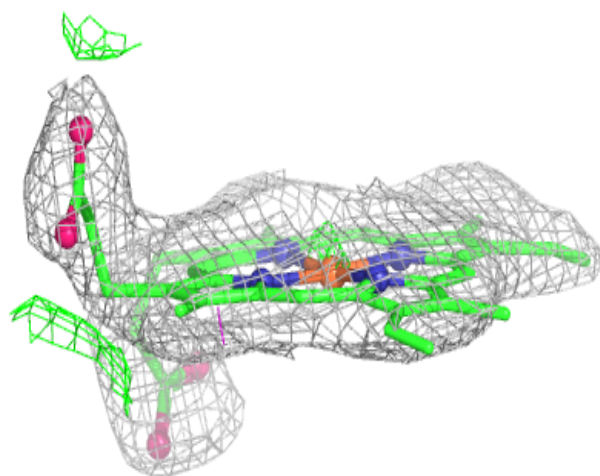
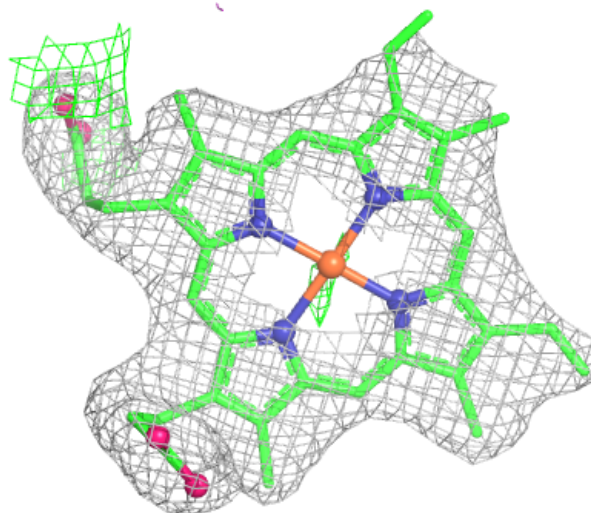
Electron density around HEM D 501:

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



Electron density around HEM B 501:

$2mF_o-DF_c$ (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.