Full wwPDB X-ray Structure Validation Report ⓘ

Jun 14, 2020 – 02:19 am BST

PDB ID	:	1DZ6
Title	:	ferrous p450cam from pseudomonas putida
Authors	:	Schlichting, I.; Berendzen, J.; Chu, K.; Stock, A.M.; Maves, S.A.; Benson, D.E.; Sweet, R.M.; Ringe, D.; Petsko, G.A.; Sligar, S.G.
Deposited on	:	2000-02-18
Resolution	:	1.90 Å(reported)

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

We welcome your comments at *validation@mail.wwpdb.org*
A user guide is available at
https://www.wwpdb.org/validation/2017/XrayValidationReportHelp
with specific help available everywhere you see the ⓘ symbol.

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

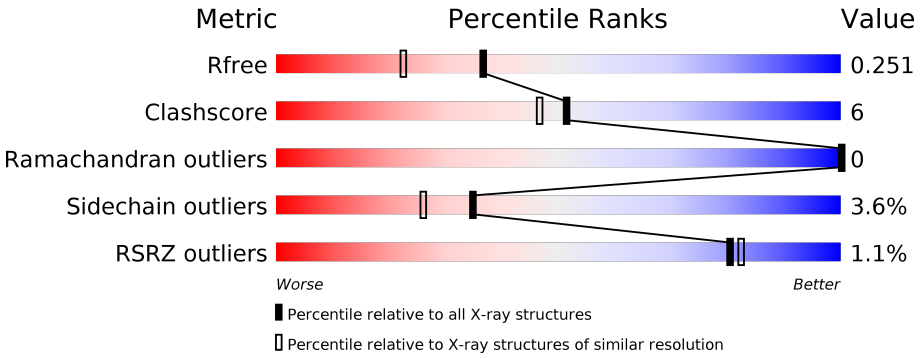
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.11
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber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

1 Overall quality at a glance (i)

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

The reported resolution of this entry is 1.90 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	6207 (1.90-1.90)
Clashscore	141614	6847 (1.90-1.90)
Ramachandran outliers	138981	6760 (1.90-1.90)
Sidechain outliers	138945	6760 (1.90-1.90)
RSRZ outliers	127900	6082 (1.90-1.90)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for >=3, 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions <=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	414	2% 77% 18%
1	B	414	80% 15%

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
4	TRS	A	503	-	X	-	-

2 Entry composition

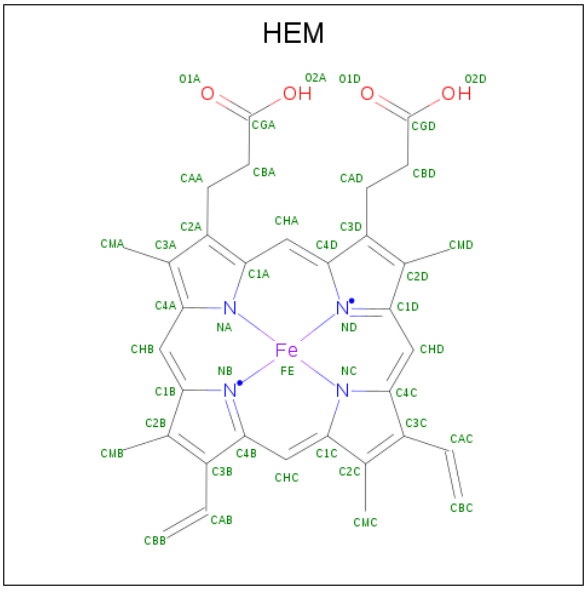
There are 6 unique types of molecules in this entry. The entry contains 7197 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 CYTOCHROME P450-CAM.

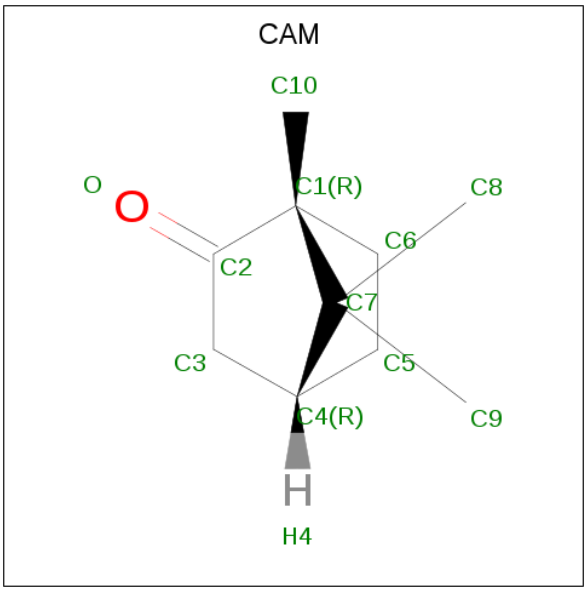
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	404	Total	C	N	O	S	0	0	0
			3200	2029	559	594	18			
1	B	405	Total	C	N	O	S	0	1	0
			3211	2035	561	597	18			

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



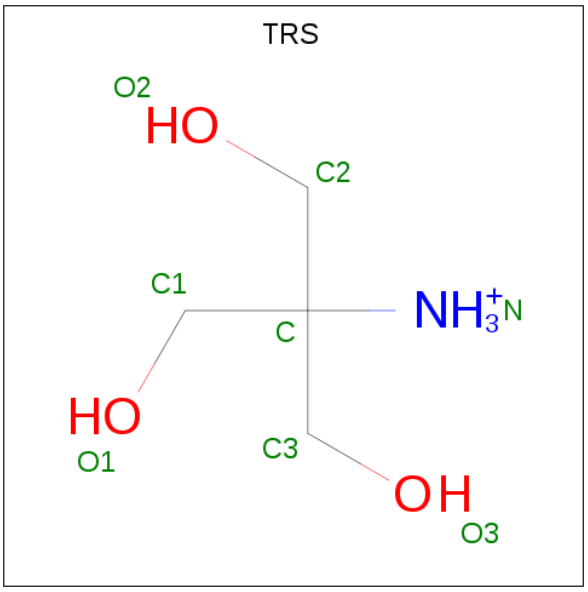
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	Fe	N	O	0	0
			43	34	1	4	4		
2	B	1	Total	C	Fe	N	O	0	0
			43	34	1	4	4		

- Molecule 3 is CAMPHOR (three-letter code: CAM) (formula: $C_{10}H_{16}O$).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total C O 11 10 1	0	0
3	B	1	Total C O 11 10 1	0	0

- Molecule 4 is 2-AMINO-2-HYDROXYMETHYL-PROPANE-1,3-DIOL (three-letter code: TRS) (formula: $C_4H_{12}NO_3$).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	1	Total C N O 8 4 1 3	0	0

- Molecule 5 is POTASSIUM ION (three-letter code: K) (formula: K).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	B	2	Total K 2 2	0	0
5	A	1	Total K 1 1	0	0

- Molecule 6 is water.

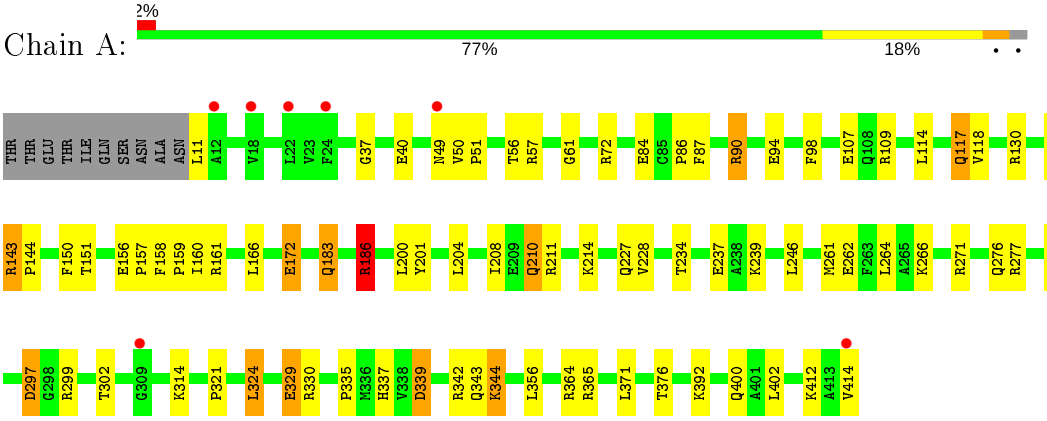
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	A	289	Total O 289 289	0	0
6	B	378	Total O 378 378	0	0

3 Residue-property plots (i)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry 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.

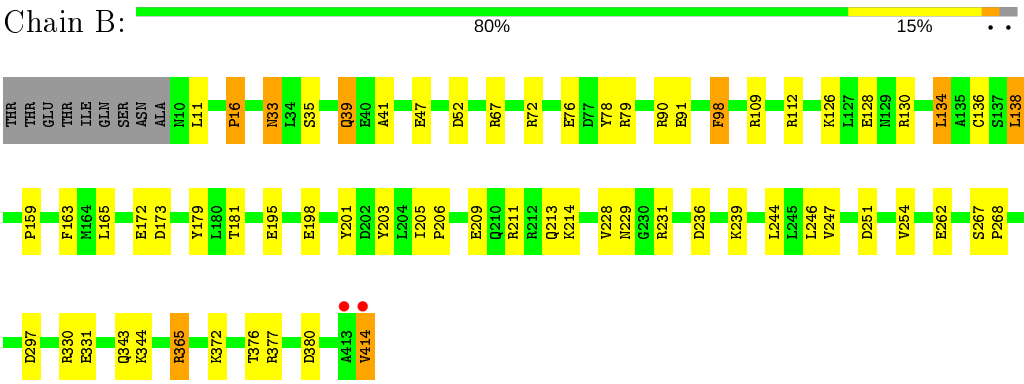
- Molecule 1: CYTOCHROME P450-CAM

Chain A:



- Molecule 1: CYTOCHROME P450-CAM

Chain B:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α, β, γ	67.80Å 62.80Å 95.50Å 90.00° 90.60° 90.00°	Depositor
Resolution (Å)	19.00 – 1.90 39.23 – 1.90	Depositor EDS
% Data completeness (in resolution range)	98.0 (19.00-1.90) 97.5 (39.23-1.90)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.09	Depositor
$< I/\sigma(I) >$ [1]	1.60 (at 1.89Å)	Xtriage
Refinement program	REFMAC	Depositor
R, R_{free}	0.188 , 0.268 0.180 , 0.251	Depositor DCC
R_{free} test set	4155 reflections (6.70%)	wwPDB-VP
Wilson B-factor (Å^2)	17.2	Xtriage
Anisotropy	0.222	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.35 , 63.5	EDS
L-test for twinning[2]	$< \|L\| > = 0.47$, $< L^2 > = 0.29$	Xtriage
Estimated twinning fraction	0.042 for h,-k,-l	Xtriage
F_o,F_c correlation	0.95	EDS
Total number of atoms	7197	wwPDB-VP
Average B, all atoms (Å^2)	23.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.60% 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

5.1 Standard geometry

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

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	0.65	0/3279	1.53	37/4454 (0.8%)
1	B	0.72	0/3294	1.59	44/4475 (1.0%)
All	All	0.69	0/6573	1.56	81/8929 (0.9%)

There are no bond length outliers.

All (81) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	B	90	ARG	NE-CZ-NH2	-13.71	113.44	120.30
1	B	280	ARG	NE-CZ-NH1	13.30	126.95	120.30
1	B	179	TYR	CB-CG-CD2	-11.27	114.24	121.00
1	A	90	ARG	NE-CZ-NH2	-11.27	114.67	120.30
1	B	179	TYR	CB-CG-CD1	10.77	127.46	121.00
1	A	186	ARG	NE-CZ-NH1	-10.73	114.93	120.30
1	A	330	ARG	NE-CZ-NH1	10.00	125.30	120.30
1	A	143	ARG	NE-CZ-NH1	9.80	125.20	120.30
1	A	271	ARG	NE-CZ-NH1	9.65	125.13	120.30
1	B	67	ARG	NE-CZ-NH1	9.60	125.10	120.30
1	B	377	ARG	NE-CZ-NH2	-9.46	115.57	120.30
1	A	90	ARG	NE-CZ-NH1	9.44	125.02	120.30
1	B	109	ARG	NE-CZ-NH2	-9.27	115.66	120.30
1	B	380	ASP	CB-CG-OD1	8.95	126.35	118.30
1	B	112	ARG	NE-CZ-NH1	8.64	124.62	120.30
1	A	365	ARG	NE-CZ-NH2	-8.57	116.01	120.30
1	B	112	ARG	NE-CZ-NH2	-8.55	116.02	120.30
1	B	280	ARG	NE-CZ-NH2	-8.48	116.06	120.30
1	A	109	ARG	NE-CZ-NH2	8.47	124.54	120.30
1	B	52	ASP	CB-CG-OD1	8.41	125.87	118.30
1	A	172	GLU	CB-CA-C	8.30	127.00	110.40
1	A	365	ARG	NE-CZ-NH1	7.89	124.25	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	B	90	ARG	CD-NE-CZ	7.89	134.65	123.60
1	A	339	ASP	CB-CG-OD1	7.56	125.11	118.30
1	B	154	TYR	CB-CG-CD1	7.54	125.53	121.00
1	B	271	ARG	CD-NE-CZ	7.42	133.99	123.60
1	B	330	ARG	NE-CZ-NH2	7.15	123.88	120.30
1	A	297	ASP	CB-CG-OD2	-7.08	111.93	118.30
1	A	342	ARG	NE-CZ-NH2	6.93	123.77	120.30
1	B	279	GLU	OE1-CD-OE2	-6.92	115.00	123.30
1	A	90	ARG	CD-NE-CZ	6.83	133.17	123.60
1	A	211	ARG	NE-CZ-NH2	6.80	123.70	120.30
1	B	251	ASP	CB-CG-OD1	6.63	124.27	118.30
1	A	277	ARG	NE-CZ-NH1	-6.62	116.99	120.30
1	A	161	ARG	NE-CZ-NH2	-6.61	117.00	120.30
1	A	271	ARG	NE-CZ-NH2	-6.53	117.03	120.30
1	B	79	ARG	NE-CZ-NH2	-6.53	117.04	120.30
1	B	195	GLU	OE1-CD-OE2	-6.44	115.57	123.30
1	B	331	GLU	OE1-CD-OE2	-6.21	115.84	123.30
1	A	98	PHE	CB-CG-CD2	-6.21	116.45	120.80
1	A	87	PHE	CB-CG-CD1	6.17	125.12	120.80
1	B	98	PHE	CB-CG-CD1	6.16	125.11	120.80
1	A	183	GLN	CA-CB-CG	6.08	126.77	113.40
1	A	291	ARG	NE-CZ-NH1	6.02	123.31	120.30
1	B	154	TYR	CB-CG-CD2	-5.96	117.43	121.00
1	A	72	ARG	NE-CZ-NH2	5.92	123.26	120.30
1	B	35	SER	O-C-N	-5.92	113.23	122.70
1	B	173	ASP	CB-CG-OD2	-5.90	112.99	118.30
1	B	72	ARG	NE-CZ-NH2	5.87	123.23	120.30
1	B	47	GLU	OE1-CD-OE2	-5.81	116.33	123.30
1	A	299	ARG	CG-CD-NE	5.80	123.98	111.80
1	A	344	LYS	N-CA-CB	5.80	121.03	110.60
1	B	330	ARG	NE-CZ-NH1	-5.74	117.43	120.30
1	A	130	ARG	NE-CZ-NH2	-5.68	117.46	120.30
1	A	330	ARG	CD-NE-CZ	5.65	131.51	123.60
1	B	76	GLU	OE1-CD-OE2	-5.64	116.53	123.30
1	A	277	ARG	NE-CZ-NH2	-5.63	117.48	120.30
1	A	277	ARG	NH1-CZ-NH2	5.50	125.45	119.40
1	B	231	ARG	CD-NE-CZ	5.45	131.23	123.60
1	B	198	GLU	OE1-CD-OE2	-5.43	116.78	123.30
1	A	98	PHE	CB-CG-CD1	5.43	124.60	120.80
1	B	203	TYR	CB-CG-CD1	-5.39	117.77	121.00
1	B	297	ASP	N-CA-CB	5.34	120.20	110.60
1	B	271	ARG	NE-CZ-NH2	-5.30	117.65	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	B	271	ARG	NE-CZ-NH1	5.30	122.95	120.30
1	B	16	PRO	O-C-N	-5.27	114.26	122.70
1	B	79	ARG	NH1-CZ-NH2	5.24	125.16	119.40
1	A	117	GLN	CA-CB-CG	5.23	124.91	113.40
1	A	166	LEU	CA-CB-CG	5.22	127.31	115.30
1	A	109	ARG	CD-NE-CZ	-5.18	116.35	123.60
1	B	78	TYR	CB-CG-CD2	-5.16	117.90	121.00
1	B	67	ARG	CG-CD-NE	5.16	122.64	111.80
1	B	78	TYR	CB-CG-CD1	5.14	124.08	121.00
1	B	262	GLU	OE1-CD-OE2	-5.11	117.16	123.30
1	B	211	ARG	CA-CB-CG	-5.06	102.27	113.40
1	A	290	ARG	NE-CZ-NH2	5.04	122.82	120.30
1	A	402	LEU	CB-CG-CD2	-5.02	102.47	111.00
1	B	79	ARG	NE-CZ-NH1	-5.01	117.79	120.30
1	A	364	ARG	NE-CZ-NH2	-5.01	117.80	120.30
1	A	210	GLN	CB-CG-CD	5.00	124.61	111.60
1	B	136	CYS	CA-CB-SG	5.00	123.01	114.00

There are no chirality outliers.

There are no planarity outliers.

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	3200	0	3152	41	0
1	B	3211	0	3161	29	0
2	A	43	0	30	3	0
2	B	43	0	30	2	0
3	A	11	0	16	0	0
3	B	11	0	16	0	0
4	A	8	0	12	0	0
5	A	1	0	0	0	0
5	B	2	0	0	0	0
6	A	289	0	0	4	0
6	B	378	0	0	2	0
All	All	7197	0	6417	72	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 6.

All (72) 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:B:236:ASP:HB3	6:B:899:HOH:O	1.74	0.87
1:A:376:THR:HG22	1:A:414:VAL:HG21	1.57	0.87
1:B:130:ARG:HD2	1:B:165:LEU:HD21	1.61	0.83
1:A:392:LYS:HD3	1:A:400:GLN:HG2	1.62	0.81
1:B:134:LEU:HD22	1:B:138:LEU:HD22	1.66	0.77
1:A:210:GLN:HE22	1:A:214:LYS:HE3	1.50	0.76
1:A:90:ARG:O	1:A:94:GLU:HG3	1.87	0.74
1:B:128:GLU:HG3	1:B:365:ARG:HH12	1.54	0.72
2:B:501:HEM:HBB2	2:B:501:HEM:HMB2	1.73	0.69
1:A:276:GLN:HE21	1:B:172:GLU:N	1.94	0.65
2:A:501:HEM:HMB2	2:A:501:HEM:HBB2	1.80	0.63
1:A:11:LEU:HD21	1:A:57:ARG:HH21	1.64	0.63
1:A:56:THR:O	1:A:61:GLY:HA2	1.99	0.63
1:B:159:PRO:HG2	1:B:254:VAL:HG22	1.79	0.62
1:A:276:GLN:HE21	1:B:172:GLU:H	1.49	0.61
1:B:376:THR:HG21	1:B:414:VAL:HG11	1.84	0.59
1:A:200:LEU:HD21	1:A:246:LEU:HD13	1.85	0.59
1:A:151:THR:HG22	6:A:732:HOH:O	2.04	0.58
1:B:372:LYS:NZ	1:B:372:LYS:HB3	2.22	0.55
1:A:114:LEU:HD13	1:A:227:GLN:O	2.06	0.55
1:B:163:PHE:HE2	1:B:246:LEU:HD23	1.72	0.54
1:B:281:ILE:HD12	1:B:372:LYS:HG2	1.88	0.54
1:A:329:GLU:HG3	1:A:335:PRO:HD3	1.90	0.53
2:B:501:HEM:HBB2	2:B:501:HEM:CMB	2.37	0.53
1:A:302:THR:O	1:A:314:LYS:HD2	2.11	0.51
1:B:33:ASN:HB3	1:B:41:ALA:HA	1.93	0.51
1:B:128:GLU:CG	1:B:365:ARG:HH12	2.22	0.50
1:A:376:THR:CG2	1:A:414:VAL:HG21	2.34	0.50
1:B:91:GLU:HB3	6:B:642:HOH:O	2.10	0.50
1:B:201:TYR:CD2	1:B:239:LYS:HE3	2.47	0.49
1:A:412:LYS:HE2	1:A:414:VAL:HG22	1.94	0.49
1:B:134:LEU:HD22	1:B:138:LEU:CD2	2.37	0.49
1:A:356:LEU:HB2	6:A:792:HOH:O	2.13	0.48
2:A:501:HEM:CMB	2:A:501:HEM:HBB2	2.43	0.48
1:A:412:LYS:HE2	1:A:414:VAL:CG2	2.44	0.47
1:A:294:LEU:HD23	1:A:294:LEU:H	1.80	0.47
1:A:262:GLU:O	1:A:266:LYS:HG3	2.15	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:151:THR:HG23	6:A:725:HOH:O	2.14	0.46
1:A:204:LEU:HD11	1:A:246:LEU:CD1	2.46	0.46
1:B:98:PHE:HB3	1:B:244:LEU:HB2	1.98	0.46
1:A:321:PRO:HB2	1:A:324:LEU:HD22	1.97	0.45
1:A:297:ASP:OD2	2:A:501:HEM:O2A	2.34	0.45
1:B:143:ARG:HB3	1:B:144:PRO:CD	2.47	0.45
1:A:210:GLN:NE2	1:A:214:LYS:HE3	2.26	0.44
1:B:209:GLU:OE2	1:B:213:GLN:NE2	2.50	0.44
1:A:50:VAL:HA	1:A:51:PRO:HD3	1.94	0.44
1:A:158:PHE:HB3	1:A:159:PRO:HD3	2.00	0.44
1:A:264:LEU:HD11	1:A:371:LEU:HD22	2.00	0.44
1:A:186:ARG:HD2	1:A:392:LYS:HG3	1.99	0.44
1:A:294:LEU:HD23	1:A:294:LEU:N	2.33	0.43
1:B:228:VAL:O	1:B:229:ASN:C	2.56	0.43
1:B:126:LYS:HD3	1:B:130:ARG:HH22	1.83	0.43
1:B:344:LYS:HA	1:B:344:LYS:HD3	1.79	0.43
1:A:107:GLU:HB2	6:A:738:HOH:O	2.17	0.43
1:B:143:ARG:N	1:B:144:PRO:HD2	2.34	0.43
1:B:39:GLN:NE2	1:B:39:GLN:H	2.15	0.43
1:A:201:TYR:HB3	1:A:239:LYS:HD2	2.00	0.43
1:A:37:GLY:HA3	1:A:40:GLU:OE1	2.18	0.43
1:A:329:GLU:HG3	1:A:335:PRO:CD	2.49	0.42
1:B:267:SER:HA	1:B:268:PRO:HD2	1.76	0.42
1:A:156:GLU:O	1:A:160:ILE:HG13	2.18	0.42
1:B:372:LYS:HB3	1:B:372:LYS:HZ3	1.83	0.42
1:A:321:PRO:HG2	1:A:324:LEU:HD22	2.01	0.42
1:B:205:ILE:N	1:B:206:PRO:HD2	2.35	0.42
1:A:234:THR:OG1	1:A:237:GLU:HG3	2.20	0.41
1:A:337:HIS:CE1	1:A:339:ASP:HB2	2.55	0.41
1:A:204:LEU:O	1:A:208:ILE:HG13	2.21	0.41
1:B:163:PHE:CE2	1:B:246:LEU:HD23	2.53	0.41
1:B:181[B]:THR:HG23	1:B:247:VAL:HG13	2.02	0.41
1:A:156:GLU:HB2	1:A:157:PRO:HD3	2.03	0.41
1:A:150:PHE:CZ	1:A:261:MET:HG3	2.56	0.41
1:A:143:ARG:HB3	1:A:144:PRO:HD3	2.03	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	402/414 (97%)	386 (96%)	16 (4%)	0	100	100
1	B	404/414 (98%)	393 (97%)	11 (3%)	0	100	100
All	All	806/828 (97%)	779 (97%)	27 (3%)	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	349/358 (98%)	334 (96%)	15 (4%)	29	19
1	B	351/358 (98%)	341 (97%)	10 (3%)	43	36
All	All	700/716 (98%)	675 (96%)	25 (4%)	35	26

All (25) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	49	ASN
1	A	84	GLU
1	A	86	PRO
1	A	117	GLN
1	A	118	VAL
1	A	137	SER
1	A	141	SER
1	A	172	GLU

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Mol	Chain	Res	Type
1	A	183	GLN
1	A	186	ARG
1	A	228	VAL
1	A	324	LEU
1	A	329	GLU
1	A	343	GLN
1	A	344	LYS
1	B	11	LEU
1	B	16	PRO
1	B	33	ASN
1	B	39	GLN
1	B	134	LEU
1	B	138	LEU
1	B	214	LYS
1	B	343	GLN
1	B	365	ARG
1	B	414	VAL

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (18) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	39	GLN
1	A	46	GLN
1	A	49	ASN
1	A	117	GLN
1	A	210	GLN
1	A	225	ASN
1	A	272	GLN
1	A	276	GLN
1	A	311	GLN
1	A	337	HIS
1	A	388	GLN
1	B	30	ASN
1	B	33	ASN
1	B	39	GLN
1	B	46	GLN
1	B	69	GLN
1	B	132	GLN
1	B	225	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 carbohydrates in this entry.

5.6 Ligand geometry

Of 8 ligands modelled in this entry, 3 are monoatomic - leaving 5 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	$\#\|Z\| > 2$
4	TRS	A	503	-	7,7,7	1.92	3 (42%)	9,9,9	2.31	4 (44%)
2	HEM	B	501	1	27,50,50	2.08	5 (18%)	17,82,82	3.06	11 (64%)
2	HEM	A	501	1	27,50,50	2.12	7 (25%)	17,82,82	2.99	8 (47%)
3	CAM	B	502	-	12,12,12	0.97	1 (8%)	20,21,21	2.46	7 (35%)
3	CAM	A	502	-	12,12,12	0.96	1 (8%)	20,21,21	2.93	10 (50%)

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
4	TRS	A	503	-	-	4/9/9/9	-
2	HEM	B	501	1	-	0/6/54/54	-
2	HEM	A	501	1	-	0/6/54/54	-
3	CAM	B	502	-	-	-	0/3/2/2

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	CAM	A	502	-	-	-	0/3/2/2

All (17) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	501	HEM	C3C-C2C	-5.60	1.32	1.40
2	A	501	HEM	C3B-C2B	-4.84	1.33	1.40
2	B	501	HEM	C3C-C2C	-4.53	1.34	1.40
2	B	501	HEM	C3B-C2B	-4.45	1.34	1.40
2	B	501	HEM	CAA-C2A	4.14	1.58	1.52
2	A	501	HEM	C3B-CAB	3.73	1.55	1.47
2	B	501	HEM	C3C-CAC	3.67	1.55	1.47
4	A	503	TRS	O3-C3	3.19	1.52	1.42
2	B	501	HEM	C3B-CAB	2.98	1.54	1.47
4	A	503	TRS	O1-C1	2.86	1.51	1.42
2	A	501	HEM	C3C-CAC	2.73	1.53	1.47
2	A	501	HEM	CAA-C2A	2.72	1.56	1.52
3	A	502	CAM	C1-C2	2.71	1.56	1.52
4	A	503	TRS	O2-C2	2.46	1.50	1.42
3	B	502	CAM	C1-C2	2.45	1.56	1.52
2	A	501	HEM	C4B-NB	2.13	1.40	1.36
2	A	501	HEM	C1C-C2C	2.06	1.47	1.42

All (40) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	A	501	HEM	CBD-CAD-C3D	6.65	124.74	112.48
3	A	502	CAM	O-C2-C3	6.22	139.39	126.47
2	B	501	HEM	CBD-CAD-C3D	6.03	123.59	112.48
3	B	502	CAM	C3-C2-C1	-5.74	95.96	107.34
2	B	501	HEM	CMA-C3A-C4A	-5.65	119.78	128.46
3	A	502	CAM	C3-C2-C1	-5.50	96.42	107.34
3	B	502	CAM	C5-C4-C3	-5.21	92.06	106.40
2	A	501	HEM	CAA-CBA-CGA	4.89	120.88	112.67
3	A	502	CAM	C5-C4-C3	-4.83	93.09	106.40
3	A	502	CAM	C7-C1-C2	4.78	108.44	100.30
2	A	501	HEM	CMA-C3A-C4A	-4.46	121.61	128.46
2	B	501	HEM	CMD-C2D-C1D	-4.27	121.90	128.46
3	A	502	CAM	C4-C3-C2	4.04	108.94	102.03
2	A	501	HEM	CMD-C2D-C1D	-4.00	122.31	128.46
4	A	503	TRS	C3-C-N	3.87	119.54	107.98
4	A	503	TRS	C3-C-C2	-3.87	98.81	110.81

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	B	501	HEM	CMC-C2C-C3C	3.78	131.75	124.68
2	A	501	HEM	CBA-CAA-C2A	3.68	119.27	112.49
3	B	502	CAM	O-C2-C3	3.55	133.85	126.47
3	B	502	CAM	C4-C3-C2	3.54	108.07	102.03
2	B	501	HEM	CMA-C3A-C2A	3.44	131.44	124.94
3	A	502	CAM	C9-C7-C4	-3.42	105.29	113.50
2	B	501	HEM	CAA-CBA-CGA	3.28	118.18	112.67
2	B	501	HEM	CMD-C2D-C3D	2.93	130.46	124.94
2	A	501	HEM	CMC-C2C-C3C	2.75	129.83	124.68
2	B	501	HEM	CBA-CAA-C2A	2.74	117.54	112.49
4	A	503	TRS	C2-C-C1	-2.73	102.34	110.81
3	B	502	CAM	C7-C1-C2	2.72	104.93	100.30
3	B	502	CAM	C6-C1-C7	2.68	107.27	101.53
2	A	501	HEM	CMA-C3A-C2A	2.66	129.96	124.94
3	B	502	CAM	C9-C7-C8	2.64	114.28	107.62
2	A	501	HEM	CMD-C2D-C3D	2.57	129.79	124.94
2	B	501	HEM	C4A-C3A-C2A	2.56	108.78	107.00
2	B	501	HEM	CMB-C2B-C3B	2.52	129.40	124.68
4	A	503	TRS	C3-C-C1	2.36	118.14	110.81
3	A	502	CAM	C5-C6-C1	-2.28	100.19	104.79
3	A	502	CAM	C6-C1-C7	2.24	106.32	101.53
3	A	502	CAM	O-C2-C1	-2.08	122.87	125.39
2	B	501	HEM	C4C-C3C-C2C	2.07	108.34	106.90
3	A	502	CAM	C8-C7-C1	2.00	117.59	113.06

There are no chirality outliers.

All (4) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	503	TRS	C1-C-C3-O3
4	A	503	TRS	N-C-C3-O3
4	A	503	TRS	C2-C-C3-O3
4	A	503	TRS	C1-C-C2-O2

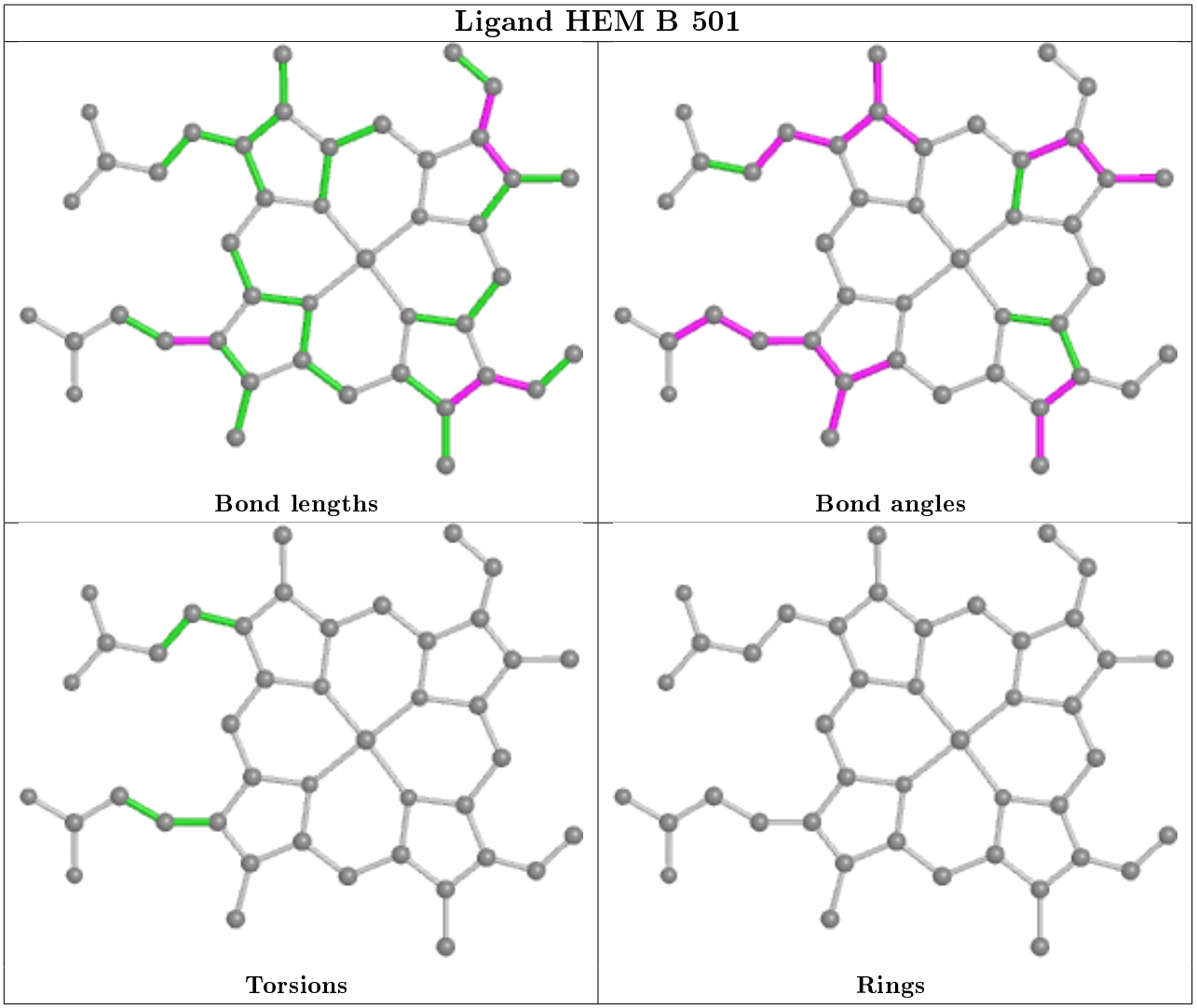
There are no ring outliers.

2 monomers are involved in 5 short contacts:

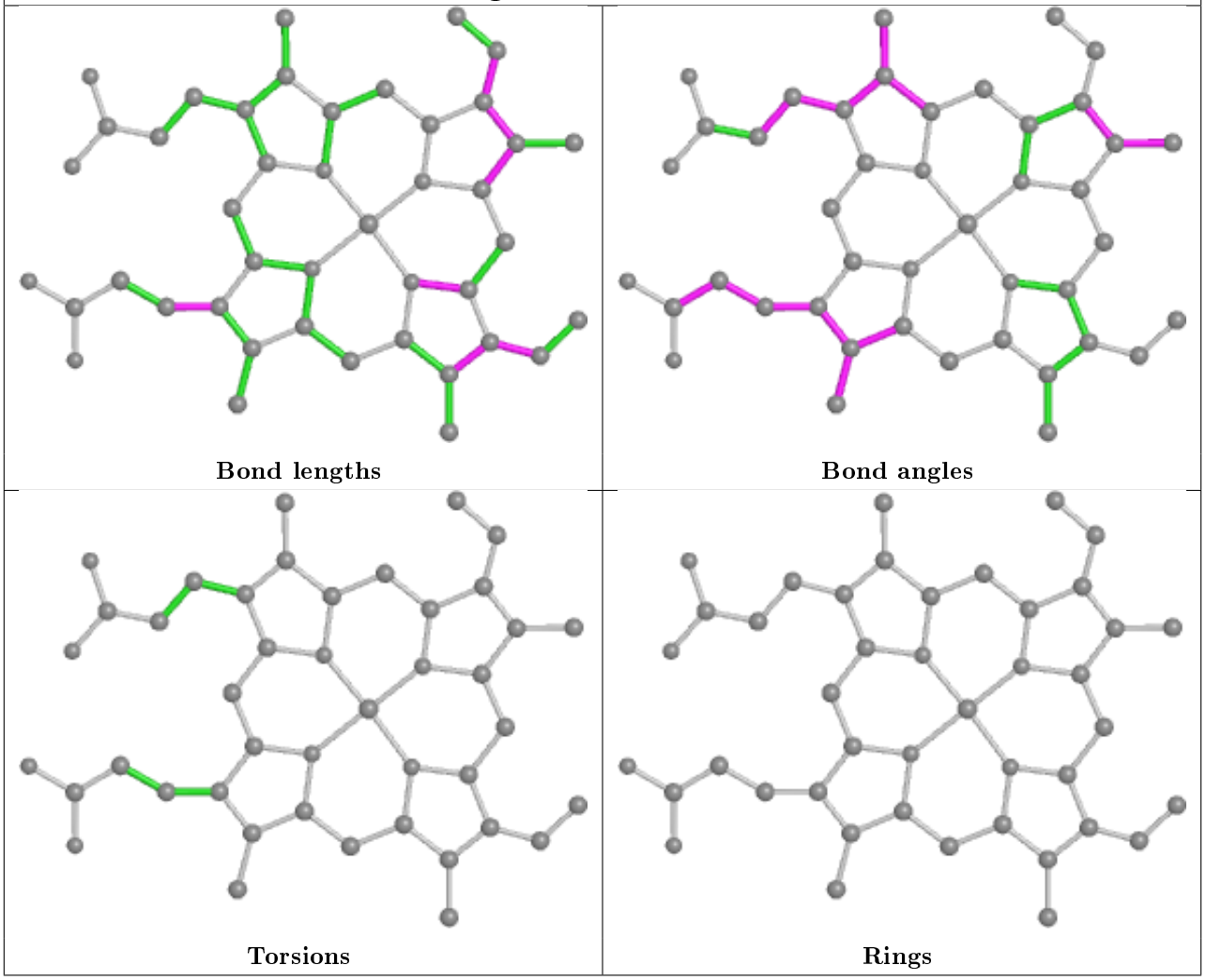
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	501	HEM	2	0
2	A	501	HEM	3	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less 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 B 501



Ligand HEM A 501



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, 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	404/414 (97%)	-0.10	7 (1%)	70	72	11, 23, 46, 64	0
1	B	405/414 (97%)	-0.34	2 (0%)	91	92	9, 18, 37, 69	0
All	All	809/828 (97%)	-0.22	9 (1%)	80	82	9, 20, 43, 69	0

All (9) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	49	ASN	3.8
1	A	22	LEU	3.0
1	B	414	VAL	3.0
1	A	24	PHE	2.7
1	A	414	VAL	2.5
1	A	18	VAL	2.4
1	A	309	GLY	2.4
1	A	12	ALA	2.1
1	B	413	ALA	2.1

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 carbohydrates 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,

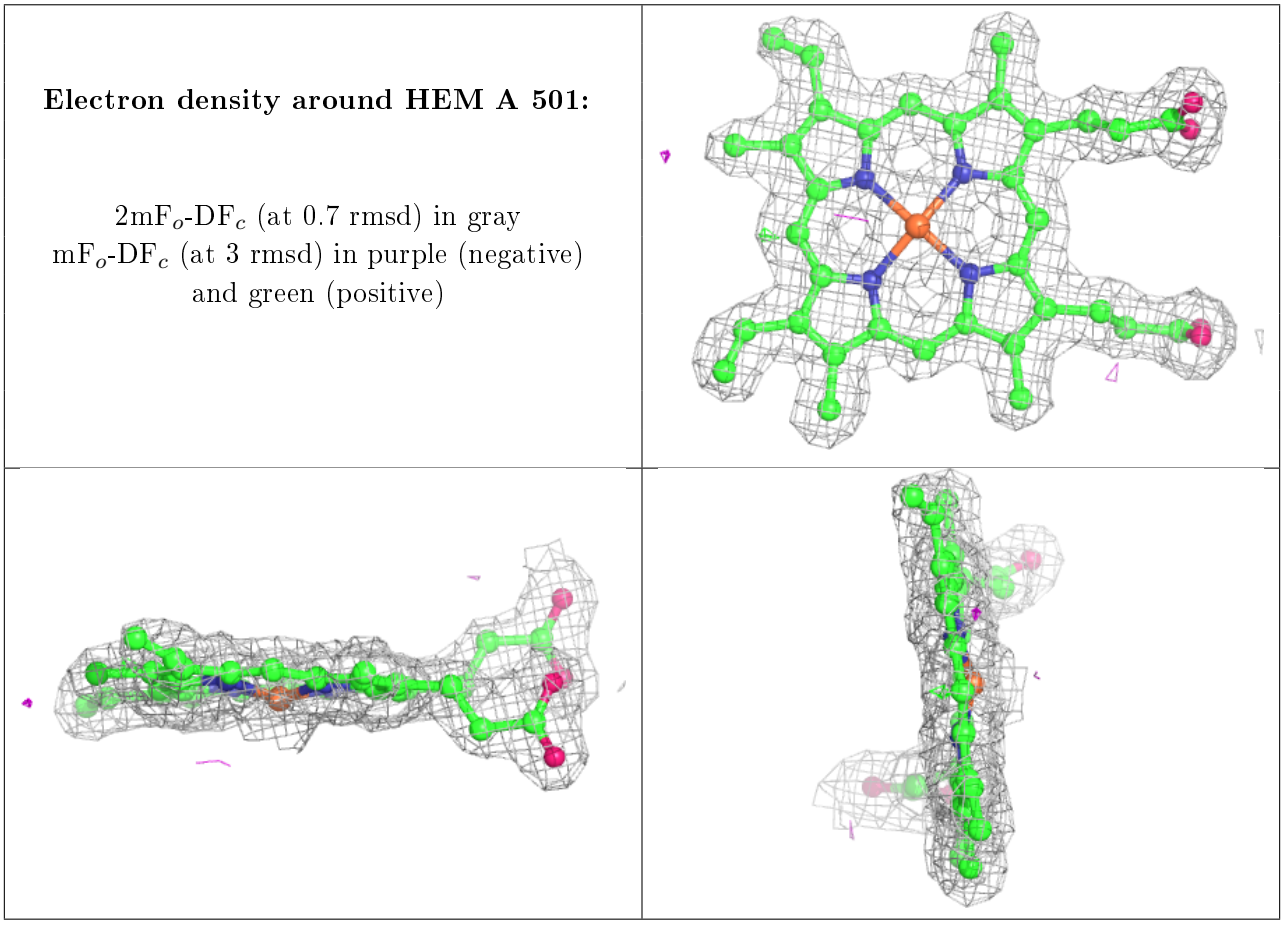
median, 95^{th} percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
3	CAM	A	502	11/11	0.93	0.15	9,14,21,21	0
4	TRS	A	503	8/8	0.95	0.07	10,17,23,27	0
3	CAM	B	502	11/11	0.96	0.19	6,11,21,23	0
2	HEM	A	501	43/43	0.98	0.13	6,17,25,31	0
2	HEM	B	501	43/43	0.98	0.12	7,13,25,33	0
5	K	A	504	1/1	0.99	0.05	21,21,21,21	0
5	K	B	504	1/1	0.99	0.03	18,18,18,18	0
5	K	B	503	1/1	1.00	0.04	12,12,12,12	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.

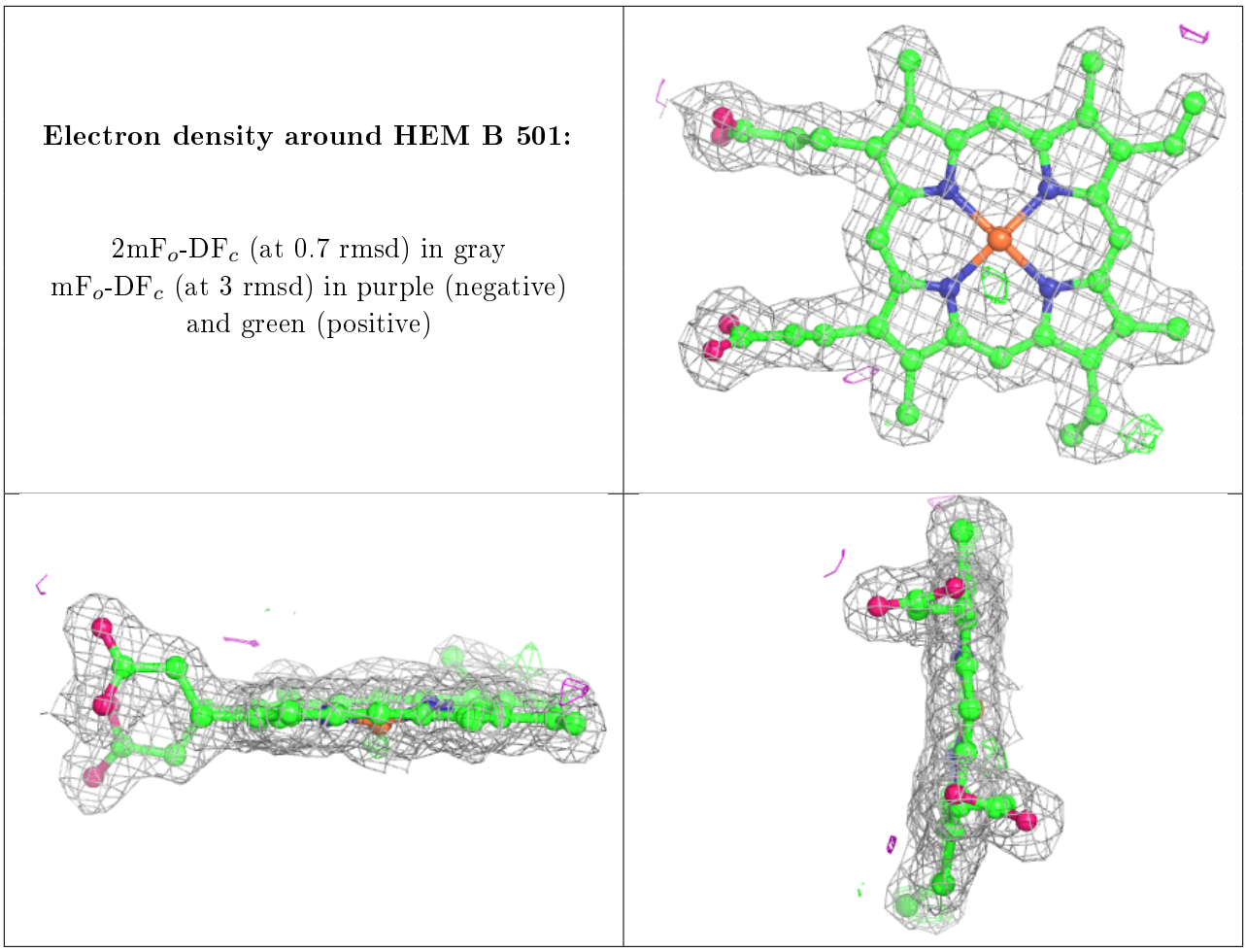
Electron density around HEM A 501:

$2mF_o$-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

There are no such residues in this entry.