



Full wwPDB X-ray Structure Validation Report ⓘ

Apr 25, 2026 – 06:55 AM EDT

PDB ID : 2GR6 / pdb_00002gr6
Title : Crystal structure of cytochrome p450cam mutant (f87w/y96f/l244a/v247l/c334a)
Authors : Rao, Z.; Wong, L.L.; Xu, F.; Bell, S.G.
Deposited on : 2006-04-22
Resolution : 2.30 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

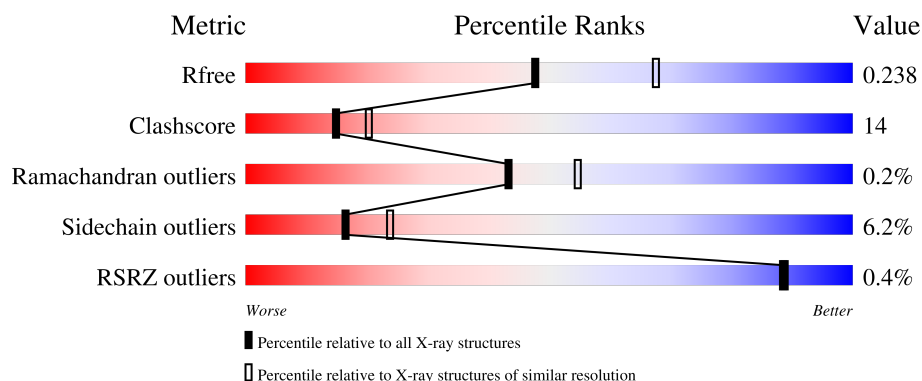
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.30 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	6319 (2.30-2.30)
Clashscore	190562	6919 (2.30-2.30)
Ramachandran outliers	187476	6854 (2.30-2.30)
Sidechain outliers	187428	6854 (2.30-2.30)
RSRZ outliers	180081	6325 (2.30-2.30)

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	405	 73% 23% .
1	B	405	 64% 34% .

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 6888 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	405	Total	C	N	O	S	0	0	0
			3206	2033	561	595	17			
1	B	405	Total	C	N	O	S	0	0	0
			3206	2033	561	595	17			

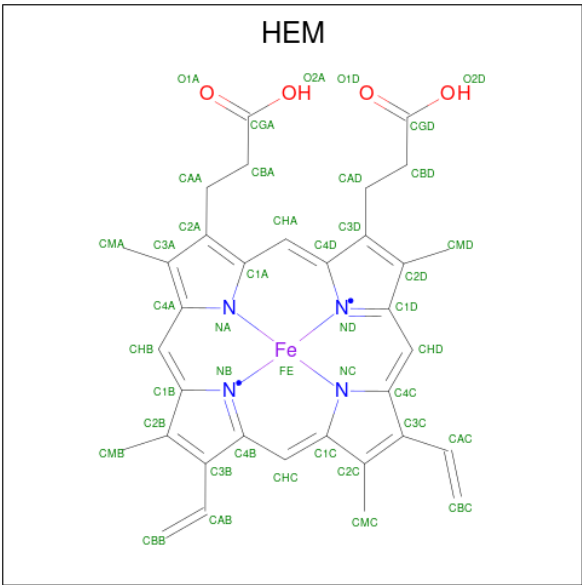
There are 10 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	87	TRP	PHE	engineered mutation	UNP P00183
A	96	PHE	TYR	engineered mutation	UNP P00183
A	244	ALA	LEU	engineered mutation	UNP P00183
A	247	LEU	VAL	engineered mutation	UNP P00183
A	334	ALA	CYS	engineered mutation	UNP P00183
B	87	TRP	PHE	engineered mutation	UNP P00183
B	96	PHE	TYR	engineered mutation	UNP P00183
B	244	ALA	LEU	engineered mutation	UNP P00183
B	247	LEU	VAL	engineered mutation	UNP P00183
B	334	ALA	CYS	engineered mutation	UNP P00183

- Molecule 2 is POTASSIUM ION (CCD ID: K) (formula: K).

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

- Molecule 3 is PROTOPORPHYRIN IX CONTAINING FE (CCD ID: HEM) (formula: C₃₄H₃₂FeN₄O₄).



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

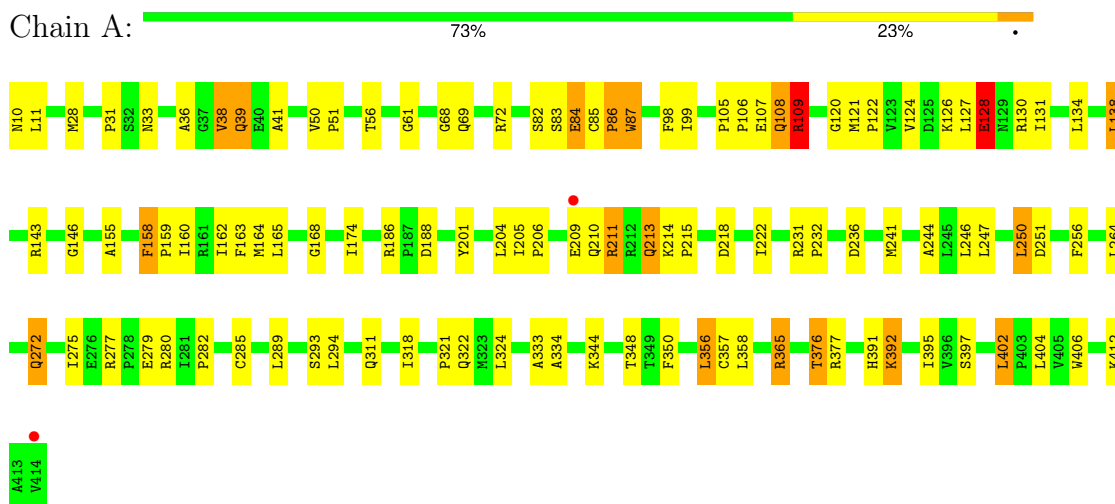
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	240	Total	O	0	0
			240	240		
4	B	148	Total	O	0	0
			148	148		

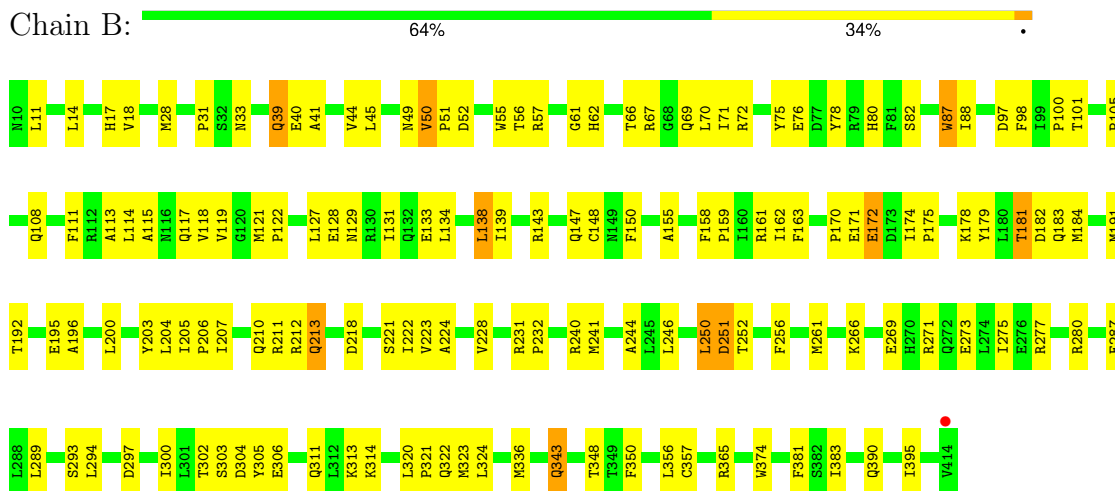
3 Residue-property plots

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: Cytochrome P450-cam



• Molecule 1: Cytochrome P450-cam



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	66.66Å 62.36Å 95.10Å 90.00° 90.37° 90.00°	Depositor
Resolution (Å)	28.00 – 2.30 28.00 – 2.30	Depositor EDS
% Data completeness (in resolution range)	(Not available) (28.00-2.30) 93.7 (28.00-2.30)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.50 (at 2.27Å)	Xtriage
Refinement program	CNS 1.0	Depositor
R, R_{free}	0.187 , 0.244 0.181 , 0.238	Depositor DCC
R_{free} test set	3605 reflections (9.97%)	wwPDB-VP
Wilson B-factor (Å ²)	32.4	Xtriage
Anisotropy	0.298	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 52.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.028 for h,-k,-l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	6888	wwPDB-VP
Average B, all atoms (Å ²)	36.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.41% 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: K, 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	0.46	1/3286 (0.0%)	0.96	14/4466 (0.3%)
1	B	0.41	0/3286	0.93	10/4466 (0.2%)
All	All	0.44	1/6572 (0.0%)	0.95	24/8932 (0.3%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	31	PRO	N-CD	9.54	1.61	1.47

All (24) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	204	LEU	N-CA-C	7.06	119.06	111.36
1	B	324	LEU	N-CA-C	6.92	118.91	111.36
1	B	204	LEU	N-CA-C	6.85	118.83	111.36
1	A	105	PRO	N-CA-C	-6.72	102.50	110.70
1	B	50	VAL	CA-C-N	6.33	126.24	119.85
1	B	50	VAL	C-N-CA	6.33	126.24	119.85
1	A	324	LEU	N-CA-C	6.14	118.94	111.82
1	A	85	CYS	CA-C-N	6.09	126.98	120.47
1	A	85	CYS	C-N-CA	6.09	126.98	120.47
1	A	31	PRO	CA-N-CD	-5.90	103.73	112.00
1	A	128	GLU	N-CA-C	5.87	118.18	111.02
1	A	109	ARG	N-CA-C	5.59	117.82	111.11
1	B	251	ASP	N-CA-C	5.47	120.20	112.93
1	B	277	ARG	CA-C-N	5.41	125.23	119.28
1	B	277	ARG	C-N-CA	5.41	125.23	119.28
1	A	158	PHE	CA-C-N	-5.28	113.59	119.19
1	A	158	PHE	C-N-CA	-5.28	113.59	119.19
1	A	214	LYS	CA-C-N	5.21	126.35	119.84

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	214	LYS	C-N-CA	5.21	126.35	119.84
1	A	174	ILE	CB-CA-C	-5.17	108.79	113.70
1	B	17	HIS	N-CA-C	5.13	119.30	113.19
1	A	31	PRO	N-CA-C	-5.10	101.65	111.69
1	B	343	GLN	N-CA-C	5.06	116.49	111.07
1	B	293	SER	N-CA-C	-5.03	103.50	110.35

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	3206	0	3153	78	0
1	B	3206	0	3153	99	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	43	0	30	1	0
3	B	43	0	30	3	0
4	A	240	0	0	3	0
4	B	148	0	0	5	0
All	All	6888	0	6366	178	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 14.

All (178) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:130:ARG:HE	1:A:165:LEU:HD21	1.09	1.13
1:B:223:VAL:HG13	1:B:241:MET:HE2	1.37	1.01
1:B:303:SER:HA	1:B:314:LYS:HG3	1.42	1.00
1:A:277:ARG:HG3	1:A:279:GLU:OE1	1.63	0.99
1:B:118:VAL:HG13	1:B:119:VAL:HG13	1.50	0.90
1:B:191:MET:HE2	1:B:196:ALA:HA	1.54	0.89

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:131:ILE:HG12	1:B:162:ILE:HD13	1.55	0.88
1:A:128:GLU:HG3	1:A:365:ARG:HE	1.39	0.87
1:B:303:SER:HA	1:B:314:LYS:CG	2.14	0.78
1:B:97:ASP:O	1:B:240:ARG:HD2	1.85	0.76
1:A:130:ARG:NE	1:A:165:LEU:HD21	1.95	0.75
1:A:128:GLU:HG3	1:A:365:ARG:NE	2.02	0.73
1:B:158:PHE:HB3	1:B:159:PRO:HD3	1.69	0.73
1:A:128:GLU:CG	1:A:365:ARG:HE	2.01	0.73
1:A:39:GLN:NE2	1:A:39:GLN:H	1.88	0.72
1:A:28:MET:HE1	1:A:395:ILE:HD13	1.72	0.71
1:B:56:THR:O	1:B:61:GLY:HA2	1.93	0.69
1:B:114:LEU:HG	1:B:241:MET:HE1	1.74	0.68
1:B:28:MET:HE1	1:B:395:ILE:HD13	1.75	0.68
1:A:294:LEU:HD23	1:A:294:LEU:H	1.60	0.67
1:A:107:GLU:H	1:A:107:GLU:CD	2.03	0.66
1:A:218:ASP:O	1:A:222:ILE:HG12	1.96	0.66
1:B:115:ALA:O	1:B:118:VAL:HG12	1.96	0.65
1:A:376:THR:O	1:A:412:LYS:HE3	1.97	0.65
1:A:205:ILE:HB	1:A:206:PRO:HD3	1.80	0.64
1:B:163:PHE:HE2	1:B:246:LEU:HD23	1.60	0.64
1:A:365:ARG:HB2	1:A:365:ARG:HH11	1.63	0.64
1:A:68:GLY:O	1:A:72:ARG:HG3	1.98	0.63
1:B:223:VAL:CG1	1:B:241:MET:HE2	2.21	0.63
1:B:181:THR:HA	1:B:184:MET:HE3	1.81	0.63
1:B:266:LYS:HD2	1:B:383:ILE:HD12	1.80	0.62
1:B:181:THR:CG2	1:B:251:ASP:HB2	2.30	0.62
1:B:131:ILE:HA	1:B:162:ILE:HD11	1.81	0.62
1:B:163:PHE:CE2	1:B:246:LEU:HD23	2.35	0.62
1:B:191:MET:HE2	1:B:196:ALA:CA	2.30	0.62
1:B:261:MET:HE3	1:B:381:PHE:CZ	2.34	0.61
1:B:294:LEU:H	1:B:294:LEU:HD23	1.66	0.61
1:A:350:PHE:HB3	1:A:357:CYS:HB3	1.82	0.61
1:B:228:VAL:HG23	1:B:228:VAL:O	2.01	0.60
1:B:205:ILE:HB	1:B:206:PRO:HD3	1.84	0.59
1:B:192:THR:OG1	1:B:195:GLU:HG3	2.03	0.58
1:B:114:LEU:CG	1:B:241:MET:HE1	2.33	0.58
1:A:134:LEU:O	1:A:138:LEU:HB2	2.03	0.58
1:B:45:LEU:HB3	1:B:66:THR:HG22	1.85	0.58
1:B:212:ARG:HG3	1:B:224:ALA:HB1	1.86	0.57
1:B:11:LEU:HD12	1:B:57:ARG:HB2	1.86	0.57
1:B:72:ARG:O	1:B:76:GLU:HG3	2.03	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:179:TYR:O	1:B:183:GLN:HG2	2.03	0.57
1:A:121:MET:HB3	1:A:122:PRO:HD3	1.86	0.57
1:A:294:LEU:HD23	1:A:294:LEU:N	2.20	0.56
1:A:106:PRO:HD2	1:A:107:GLU:OE2	2.06	0.56
1:A:87:TRP:CZ2	1:A:395:ILE:HG21	2.41	0.56
1:B:266:LYS:HD2	1:B:383:ILE:CD1	2.35	0.56
1:B:306:GLU:HB2	1:B:311:GLN:NE2	2.20	0.55
1:A:146:GLY:HA2	1:A:406:TRP:CD1	2.42	0.55
1:B:170:PRO:HB2	1:B:172:GLU:HG2	1.89	0.55
1:B:87:TRP:CZ2	1:B:395:ILE:HG21	2.43	0.54
1:A:272:GLN:HA	1:A:275:ILE:HD12	1.89	0.54
1:B:114:LEU:CD2	1:B:241:MET:HE1	2.38	0.54
1:A:128:GLU:CD	1:A:365:ARG:HE	2.16	0.54
1:B:39:GLN:NE2	1:B:39:GLN:H	2.07	0.53
1:A:186:ARG:NH2	1:A:251:ASP:OD2	2.35	0.53
1:A:82:SER:OG	1:A:84:GLU:HB2	2.10	0.52
1:B:139:ILE:HD13	1:B:374:TRP:HA	1.92	0.52
1:A:376:THR:HG22	1:A:377:ARG:HG3	1.91	0.52
1:B:75:TYR:CZ	1:B:320:LEU:HB2	2.45	0.51
1:B:200:LEU:HD11	1:B:246:LEU:CD1	2.41	0.51
1:B:98:PHE:HB3	1:B:244:ALA:HB2	1.92	0.51
1:B:184:MET:HE1	1:B:200:LEU:HD22	1.91	0.51
1:A:213:GLN:C	1:A:215:PRO:HD3	2.36	0.51
1:A:56:THR:O	1:A:61:GLY:HA2	2.10	0.50
1:A:33:ASN:ND2	1:A:36:ALA:HB3	2.27	0.50
1:B:181:THR:HG21	1:B:251:ASP:HB2	1.94	0.50
1:A:127:LEU:O	1:A:131:ILE:HG13	2.12	0.50
1:B:150:PHE:CE1	1:B:261:MET:HG2	2.46	0.50
1:A:38:VAL:HG11	1:A:397:SER:HB3	1.95	0.49
1:A:277:ARG:HB3	1:A:280:ARG:HG3	1.93	0.49
1:B:150:PHE:CE2	1:B:155:ALA:HB2	2.47	0.49
1:B:200:LEU:HD11	1:B:246:LEU:HD12	1.94	0.49
1:B:62:HIS:CG	1:B:88:ILE:HD13	2.47	0.49
1:A:33:ASN:HB3	1:A:41:ALA:HA	1.95	0.49
1:A:120:GLY:O	1:A:124:VAL:HG23	2.12	0.49
1:B:207:ILE:HG23	1:B:211:ARG:NH1	2.28	0.49
1:B:256:PHE:CE2	1:B:289:LEU:HD23	2.48	0.48
1:A:38:VAL:HG22	1:A:391:HIS:HB3	1.94	0.48
1:B:171:GLU:HG3	4:B:2439:HOH:O	2.12	0.48
1:A:98:PHE:HB3	1:A:244:ALA:HB2	1.96	0.48
1:A:402:LEU:HD22	1:A:404:LEU:CD2	2.43	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:163:PHE:HE2	1:B:246:LEU:CD2	2.26	0.48
1:A:87:TRP:CH2	1:A:395:ILE:HG21	2.48	0.48
1:A:201:TYR:O	1:A:205:ILE:HG13	2.14	0.48
1:A:246:LEU:HD12	1:A:250:LEU:HD22	1.96	0.48
1:A:83:SER:O	1:A:86:PRO:HD3	2.14	0.48
1:A:158:PHE:O	1:A:162:ILE:HG12	2.13	0.48
1:B:273:GLU:OE1	1:B:280:ARG:NH2	2.46	0.48
1:A:365:ARG:HB2	1:A:365:ARG:NH1	2.29	0.47
1:A:188:ASP:OD2	1:A:188:ASP:C	2.57	0.47
1:B:174:ILE:HB	1:B:175:PRO:HD3	1.96	0.47
1:A:402:LEU:HD22	1:A:404:LEU:HD21	1.97	0.47
1:B:131:ILE:HA	1:B:162:ILE:CD1	2.45	0.47
1:A:99:ILE:O	1:A:241:MET:HA	2.15	0.46
1:B:206:PRO:O	1:B:210:GLN:HG3	2.15	0.46
1:A:168:GLY:C	1:A:211:ARG:HH22	2.23	0.46
1:A:246:LEU:O	1:A:250:LEU:HB2	2.15	0.46
1:B:41:ALA:O	1:B:44:VAL:HG22	2.16	0.46
1:B:269:GLU:HB2	4:B:2561:HOH:O	2.14	0.46
1:A:146:GLY:HA2	1:A:406:TRP:NE1	2.30	0.46
1:A:294:LEU:N	1:A:294:LEU:CD2	2.79	0.46
1:B:49:ASN:ND2	4:B:2485:HOH:O	2.49	0.46
1:B:271:ARG:O	1:B:275:ILE:HG13	2.16	0.46
1:B:287:GLU:HA	1:B:287:GLU:OE1	2.16	0.46
1:B:114:LEU:HD23	1:B:241:MET:HE1	1.97	0.46
1:B:118:VAL:CG1	1:B:119:VAL:HG13	2.35	0.46
1:A:39:GLN:H	1:A:39:GLN:HE21	1.61	0.45
1:A:213:GLN:O	1:A:215:PRO:HD3	2.17	0.45
1:A:109:ARG:HD2	4:A:1650:HOH:O	2.15	0.45
1:B:14:LEU:HD11	1:B:18:VAL:CG1	2.46	0.45
1:B:300:ILE:O	1:B:300:ILE:HG23	2.15	0.45
1:A:311:GLN:NE2	4:A:1472:HOH:O	2.43	0.45
1:B:31:PRO:HB2	1:B:41:ALA:HB1	1.98	0.45
1:A:322:GLN:HB3	1:A:348:THR:O	2.17	0.45
1:B:82:SER:HB2	1:B:302:THR:CG2	2.47	0.45
1:B:67:ARG:O	1:B:71:ILE:HG13	2.17	0.44
1:A:168:GLY:O	1:A:211:ARG:NH2	2.45	0.44
1:B:121:MET:HB3	1:B:122:PRO:HD3	1.99	0.44
1:B:147:GLN:O	1:B:148:CYS:HB3	2.18	0.44
1:A:163:PHE:CE2	1:A:246:LEU:HD13	2.52	0.44
1:B:218:ASP:O	1:B:222:ILE:HG12	2.16	0.44
1:B:322:GLN:HB3	1:B:348:THR:O	2.18	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:38:VAL:HG22	1:A:391:HIS:ND1	2.32	0.43
1:B:294:LEU:HD23	1:B:294:LEU:N	2.30	0.43
1:B:87:TRP:CH2	1:B:395:ILE:HG21	2.53	0.43
1:B:161:ARG:NH1	4:B:2428:HOH:O	2.47	0.43
1:A:277:ARG:O	1:A:280:ARG:HG3	2.20	0.42
1:B:40:GLU:HG3	1:B:336:MET:HE2	2.00	0.42
1:B:178:LYS:HE3	1:B:182:ASP:OD1	2.18	0.42
1:B:350:PHE:HB3	1:B:357:CYS:HB3	2.01	0.42
1:A:256:PHE:CE2	1:A:289:LEU:HA	2.54	0.42
1:B:231:ARG:HA	1:B:232:PRO:HD3	1.92	0.42
1:B:304:ASP:OD1	1:B:313:LYS:HD2	2.20	0.42
1:A:264:LEU:HD23	1:A:264:LEU:HA	1.89	0.42
1:B:113:ALA:O	1:B:117:GLN:HG3	2.19	0.42
1:A:186:ARG:HD2	1:A:392:LYS:HB3	2.01	0.42
1:A:231:ARG:HA	1:A:232:PRO:HD3	1.89	0.42
1:B:55:TRP:CD1	1:B:55:TRP:C	2.97	0.42
1:B:129:ASN:HB3	4:B:2424:HOH:O	2.18	0.42
1:A:50:VAL:HA	1:A:51:PRO:HD3	1.91	0.42
1:A:122:PRO:O	1:A:126:LYS:HG2	2.19	0.42
1:B:143:ARG:HG2	1:B:143:ARG:HH11	1.85	0.42
1:B:211:ARG:HG3	1:B:221:SER:OG	2.20	0.42
1:B:246:LEU:O	1:B:250:LEU:HB2	2.19	0.42
1:A:160:ILE:O	1:A:164:MET:HG2	2.19	0.41
1:A:210:GLN:OE1	4:A:1579:HOH:O	2.22	0.41
1:B:80:HIS:CD2	1:B:305:TYR:CG	3.07	0.41
1:B:50:VAL:HA	1:B:51:PRO:HD3	1.77	0.41
1:A:108:GLN:HE21	1:A:108:GLN:HB3	1.55	0.41
1:B:78:TYR:CE1	1:B:105:PRO:HG2	2.56	0.41
1:A:282:PRO:O	1:A:285:CYS:HB3	2.20	0.41
1:B:62:HIS:CD2	1:B:88:ILE:HD13	2.55	0.41
1:B:100:PRO:HB3	1:B:111:PHE:HB2	2.02	0.41
1:B:200:LEU:O	1:B:203:TYR:HB3	2.20	0.41
1:A:155:ALA:O	1:A:159:PRO:HD2	2.20	0.41
1:B:127:LEU:O	1:B:128:GLU:C	2.63	0.41
1:B:297:ASP:OD2	3:B:417:HEM:O2A	2.39	0.41
1:A:138:LEU:HD12	1:A:138:LEU:HA	1.93	0.41
1:B:52:ASP:O	1:B:70:LEU:HD22	2.21	0.41
1:B:101:THR:HA	3:B:417:HEM:O2D	2.21	0.41
1:B:213:GLN:HE21	1:B:213:GLN:HB2	1.66	0.41
1:B:223:VAL:HG22	1:B:241:MET:HE3	2.02	0.41
1:A:143:ARG:HH11	1:A:143:ARG:HG2	1.85	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:163:PHE:HE2	1:A:246:LEU:HD13	1.85	0.41
1:A:318:ILE:O	1:A:318:ILE:HG23	2.21	0.41
1:A:356:LEU:HD12	1:A:356:LEU:HA	1.91	0.40
1:A:358:LEU:HD12	3:A:417:HEM:HMD1	2.03	0.40
1:A:333:ALA:O	1:A:334:ALA:C	2.64	0.40
1:B:134:LEU:O	1:B:138:LEU:HB2	2.21	0.40
1:B:45:LEU:HB3	1:B:66:THR:CG2	2.51	0.40
3:B:417:HEM:HBB2	3:B:417:HEM:HMB1	2.02	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	403/405 (100%)	383 (95%)	19 (5%)	1 (0%)	43	55
1	B	403/405 (100%)	389 (96%)	13 (3%)	1 (0%)	43	55
All	All	806/810 (100%)	772 (96%)	32 (4%)	2 (0%)	43	55

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	321	PRO
1	B	321	PRO

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	348/348 (100%)	322 (92%)	26 (8%)	12	17
1	B	348/348 (100%)	331 (95%)	17 (5%)	22	34
All	All	696/696 (100%)	653 (94%)	43 (6%)	16	24

All (43) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	10	ASN
1	A	11	LEU
1	A	38	VAL
1	A	39	GLN
1	A	69	GLN
1	A	84	GLU
1	A	86	PRO
1	A	87	TRP
1	A	108	GLN
1	A	109	ARG
1	A	128	GLU
1	A	138	LEU
1	A	209	GLU
1	A	211	ARG
1	A	213	GLN
1	A	236	ASP
1	A	247	LEU
1	A	250	LEU
1	A	272	GLN
1	A	293	SER
1	A	344	LYS
1	A	356	LEU
1	A	365	ARG
1	A	376	THR
1	A	392	LYS
1	A	402	LEU
1	B	33	ASN
1	B	39	GLN
1	B	69	GLN
1	B	87	TRP
1	B	108	GLN
1	B	133	GLU
1	B	138	LEU
1	B	172	GLU

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Mol	Chain	Res	Type
1	B	181	THR
1	B	213	GLN
1	B	250	LEU
1	B	252	THR
1	B	323	MET
1	B	343	GLN
1	B	356	LEU
1	B	365	ARG
1	B	390	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (31) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	30	ASN
1	A	33	ASN
1	A	39	GLN
1	A	46	GLN
1	A	69	GLN
1	A	108	GLN
1	A	110	GLN
1	A	129	ASN
1	A	132	GLN
1	A	147	GLN
1	A	213	GLN
1	A	225	ASN
1	A	229	ASN
1	A	311	GLN
1	A	337	HIS
1	A	388	GLN
1	B	33	ASN
1	B	39	GLN
1	B	46	GLN
1	B	80	HIS
1	B	108	GLN
1	B	132	GLN
1	B	149	ASN
1	B	210	GLN
1	B	213	GLN
1	B	225	ASN
1	B	227	GLN
1	B	229	ASN
1	B	272	GLN

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Mol	Chain	Res	Type
1	B	317	GLN
1	B	388	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

Of 4 ligands modelled in this entry, 2 are monoatomic - leaving 2 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	$\# Z > 2$
3	HEM	A	417	4,1	50,50,50	1.97	15 (30%)	67,82,82	1.60	13 (19%)
3	HEM	B	417	4,1	50,50,50	2.00	16 (32%)	67,82,82	1.58	13 (19%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	HEM	A	417	4,1	-	4/14/54/54	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	HEM	B	417	4,1	-	4/14/54/54	-

All (31) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	417	HEM	FE-ND	4.73	2.09	1.94
3	A	417	HEM	FE-ND	4.59	2.09	1.94
3	A	417	HEM	CAC-C3C	4.59	1.59	1.47
3	B	417	HEM	CAC-C3C	4.27	1.58	1.47
3	B	417	HEM	CBD-CGD	4.22	1.60	1.50
3	A	417	HEM	CBD-CGD	4.08	1.60	1.50
3	A	417	HEM	CAB-C3B	3.85	1.57	1.47
3	B	417	HEM	CAB-C3B	3.83	1.57	1.47
3	A	417	HEM	CBA-CGA	3.61	1.58	1.50
3	B	417	HEM	FE-NB	3.56	2.05	1.94
3	A	417	HEM	FE-NB	3.36	2.05	1.94
3	A	417	HEM	CMC-C2C	3.34	1.57	1.50
3	B	417	HEM	CBA-CGA	3.34	1.58	1.50
3	B	417	HEM	FE-NC	3.27	2.06	1.95
3	B	417	HEM	CMC-C2C	2.74	1.56	1.50
3	B	417	HEM	CMD-C2D	2.73	1.56	1.50
3	A	417	HEM	FE-NC	2.68	2.04	1.95
3	A	417	HEM	CAA-C2A	2.59	1.58	1.51
3	B	417	HEM	CMB-C2B	2.53	1.56	1.50
3	A	417	HEM	CMD-C2D	2.50	1.55	1.50
3	A	417	HEM	FE-NA	2.47	2.03	1.95
3	B	417	HEM	C3B-C2B	-2.47	1.32	1.37
3	B	417	HEM	CAD-C3D	2.44	1.57	1.51
3	A	417	HEM	CMB-C2B	2.43	1.55	1.50
3	B	417	HEM	C1D-C2D	2.35	1.49	1.44
3	B	417	HEM	CAA-C2A	2.32	1.57	1.51
3	B	417	HEM	FE-NA	2.31	2.02	1.95
3	A	417	HEM	C3B-C2B	-2.22	1.32	1.37
3	A	417	HEM	CAD-C3D	2.17	1.56	1.51
3	A	417	HEM	O2D-CGD	-2.11	1.23	1.30
3	B	417	HEM	C4A-C3A	2.00	1.47	1.43

All (26) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	417	HEM	O2D-CGD-O1D	4.60	135.16	123.33
3	A	417	HEM	O2D-CGD-O1D	4.56	135.05	123.33

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	417	HEM	O1D-CGD-CBD	-4.00	110.41	123.09
3	A	417	HEM	C4D-ND-C1D	3.89	109.81	105.21
3	B	417	HEM	O2A-CGA-O1A	3.88	133.31	123.33
3	A	417	HEM	O1D-CGD-CBD	-3.85	110.89	123.09
3	B	417	HEM	C4D-ND-C1D	3.81	109.72	105.21
3	A	417	HEM	O2A-CGA-O1A	3.61	132.63	123.33
3	A	417	HEM	CHD-C1D-ND	3.46	128.14	124.42
3	A	417	HEM	C3D-C4D-ND	-3.24	106.62	110.17
3	B	417	HEM	CHD-C1D-ND	3.22	127.89	124.42
3	A	417	HEM	CHA-C4D-ND	3.02	128.10	124.37
3	B	417	HEM	C3D-C4D-ND	-2.93	106.96	110.17
3	A	417	HEM	CHB-C4A-NA	2.88	129.09	123.86
3	B	417	HEM	CHB-C4A-NA	2.86	129.05	123.86
3	B	417	HEM	CHA-C4D-ND	2.65	127.64	124.37
3	B	417	HEM	O2A-CGA-CBA	-2.45	106.26	114.00
3	A	417	HEM	CBA-CAA-C2A	2.40	119.18	112.53
3	A	417	HEM	O2A-CGA-CBA	-2.29	106.76	114.00
3	B	417	HEM	CMB-C2B-C1B	-2.19	121.61	125.03
3	B	417	HEM	CBC-CAC-C3C	-2.19	116.60	127.53
3	A	417	HEM	CBC-CAC-C3C	-2.17	116.69	127.53
3	A	417	HEM	CMA-C3A-C2A	2.15	130.19	125.62
3	B	417	HEM	C2D-C1D-ND	-2.12	107.45	109.90
3	A	417	HEM	CMB-C2B-C1B	-2.09	121.77	125.03
3	B	417	HEM	CMA-C3A-C2A	2.05	129.97	125.62

There are no chirality outliers.

All (8) torsion outliers are listed below:

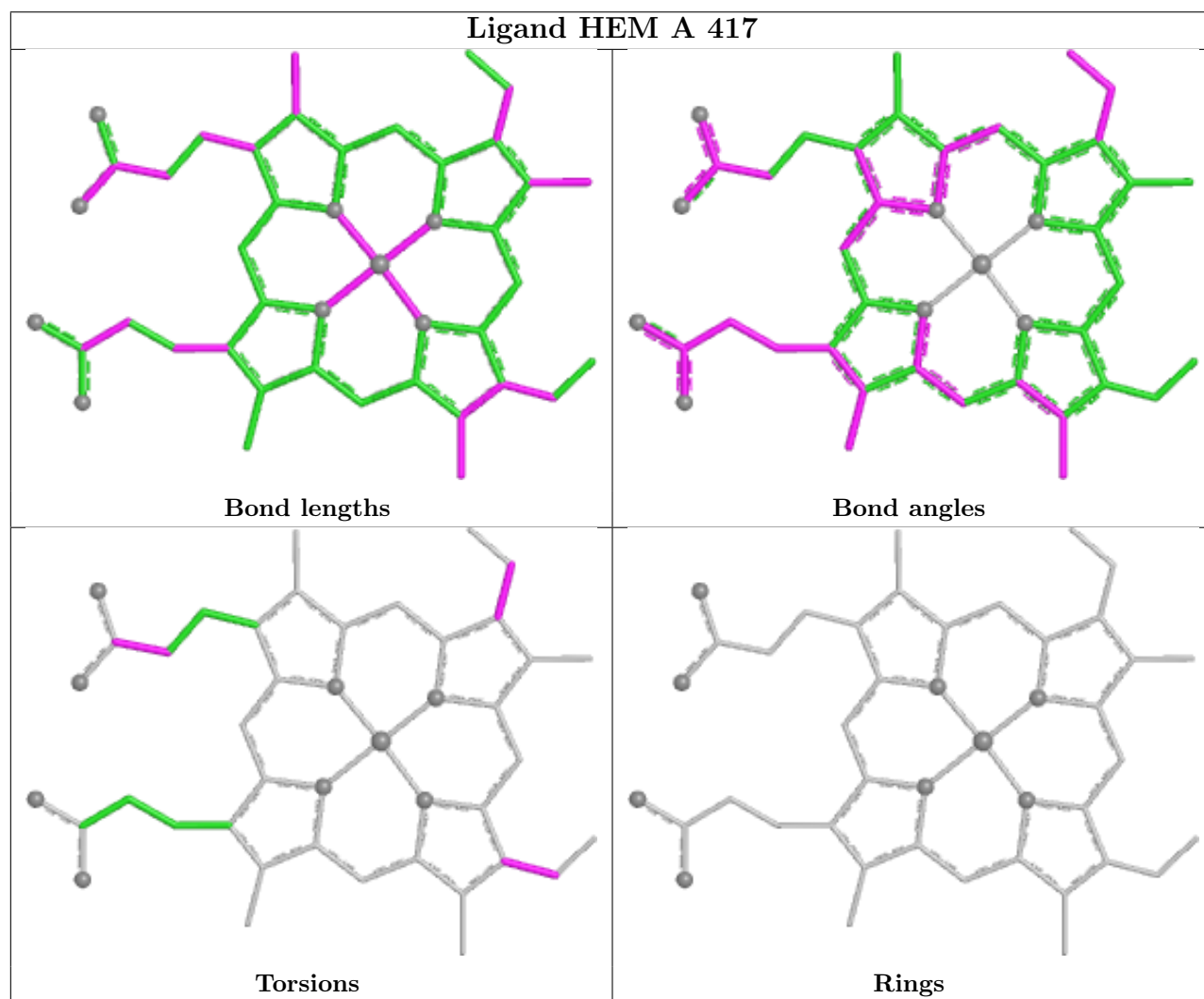
Mol	Chain	Res	Type	Atoms
3	B	417	HEM	C2C-C3C-CAC-CBC
3	B	417	HEM	C4C-C3C-CAC-CBC
3	A	417	HEM	C2C-C3C-CAC-CBC
3	A	417	HEM	C4C-C3C-CAC-CBC
3	A	417	HEM	C2B-C3B-CAB-CBB
3	B	417	HEM	CAD-CBD-CGD-O2D
3	A	417	HEM	CAD-CBD-CGD-O2D
3	B	417	HEM	CAD-CBD-CGD-O1D

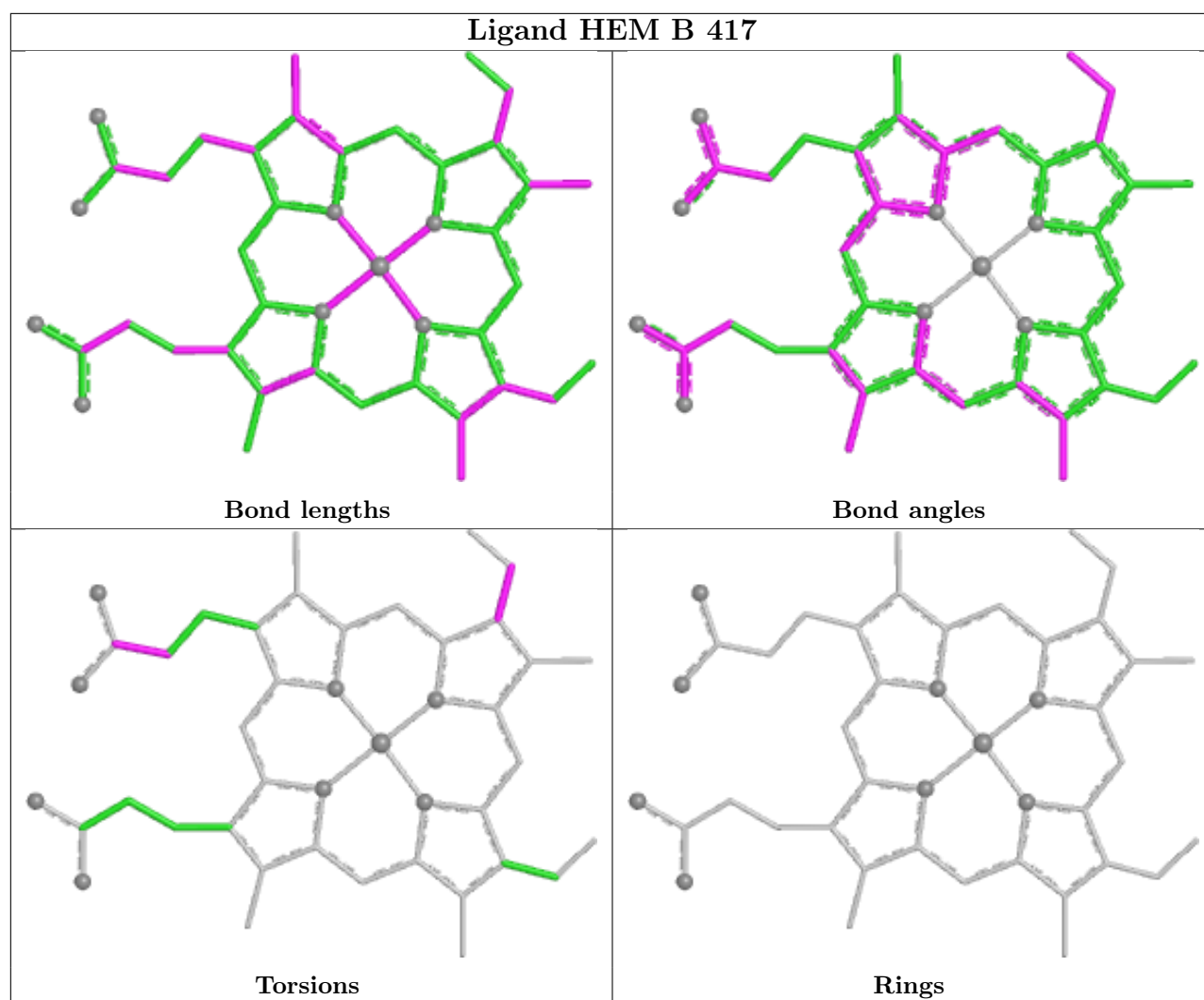
There are no ring outliers.

2 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	417	HEM	1	0
3	B	417	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 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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	405/405 (100%)	-0.31	2 (0%) 87 87	20, 31, 48, 63	0
1	B	405/405 (100%)	-0.05	1 (0%) 91 91	25, 38, 58, 73	0
All	All	810/810 (100%)	-0.18	3 (0%) 88 89	20, 35, 55, 73	0

All (3) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	414	VAL	2.4
1	A	209	GLU	2.2
1	A	414	VAL	2.1

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

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
2	K	B	2418	1/1	0.75	0.13	56,56,56,56	0
2	K	A	1418	1/1	0.96	0.04	33,33,33,33	0

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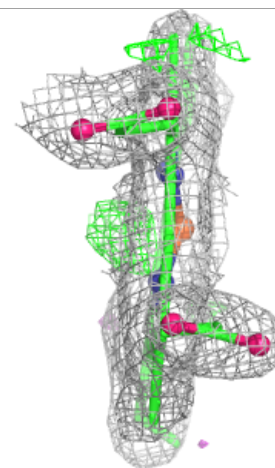
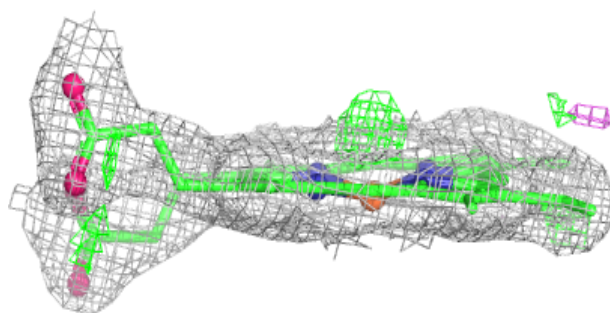
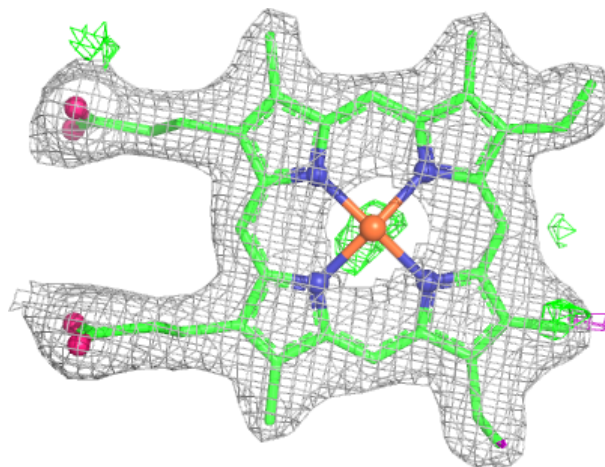
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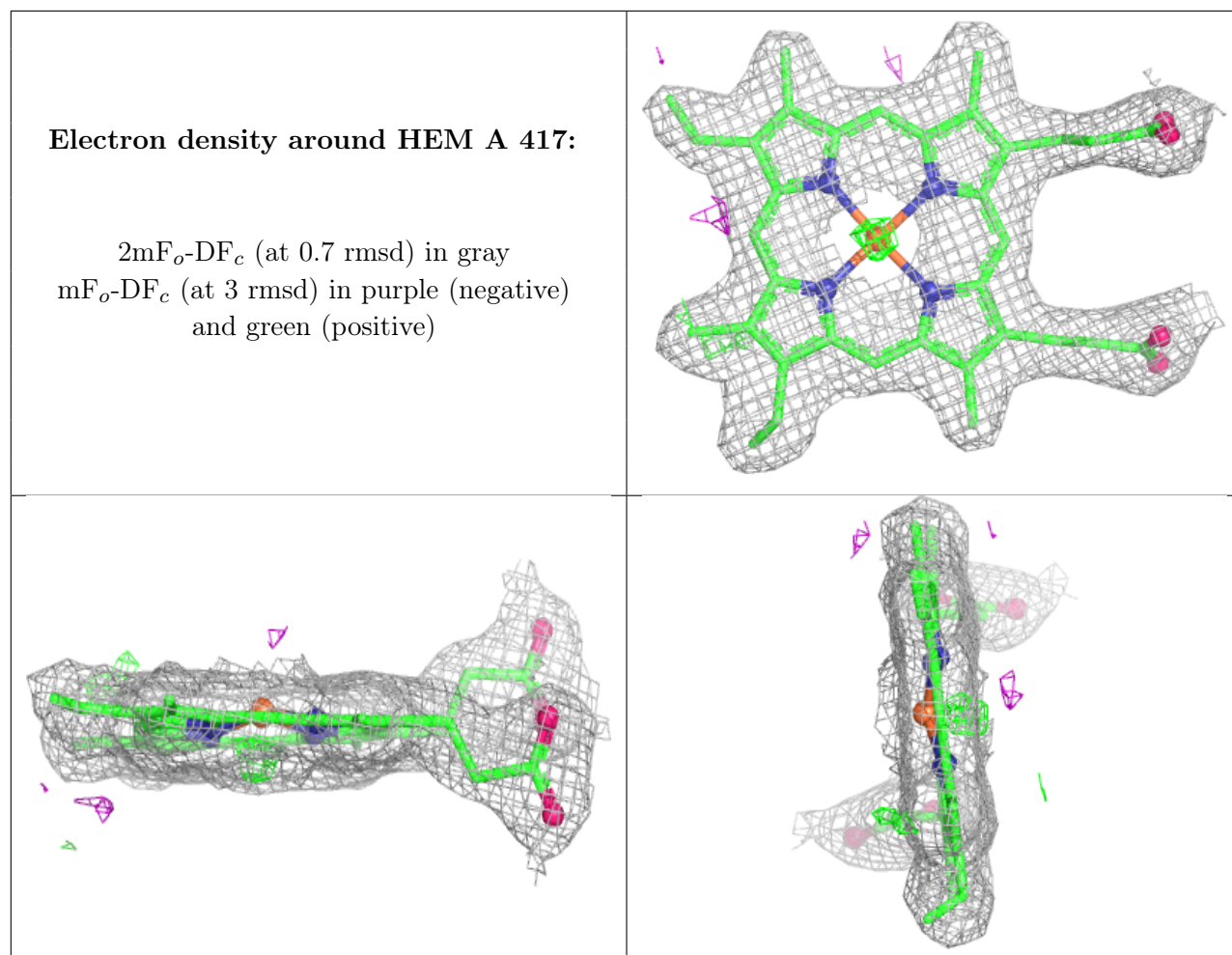
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	HEM	B	417	43/43	0.96	0.07	26,31,35,36	0
3	HEM	A	417	43/43	0.97	0.06	17,23,27,29	0

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

Electron density around HEM B 417:

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.