Full wwPDB X-ray Structure Validation Report (i)

Aug 22, 2026 – 03:48 am BST

PDB ID	:	1O76 / pdb_00001o76
Title	:	CYANIDE COMPLEX OF P450CAM FROM PSEUDOMONAS PUTIDA
Authors	:	Fedorov, R.; Ghosh, D.; Schlichting, I.
Deposited on	:	2002-10-23
Resolution	:	1.80 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 (i) 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 (i)) were used in the production of this report:

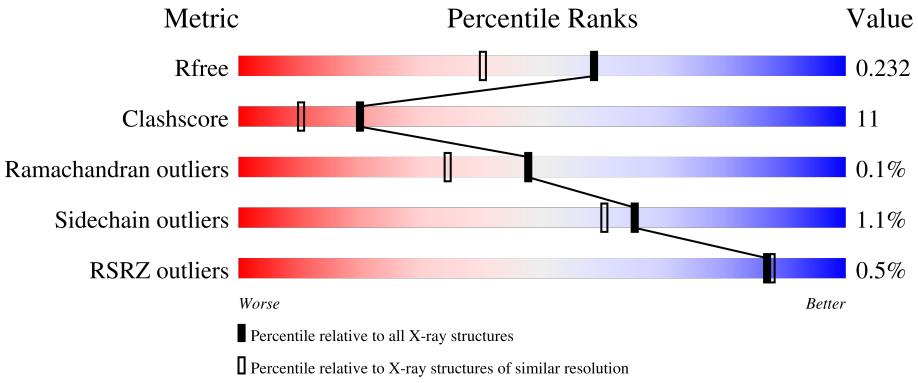
MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	1.8.4, CSD as541be (2020)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.015 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.50

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 1.80 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	7662 (1.80-1.80)
Clashscore	190562	8479 (1.80-1.80)
Ramachandran outliers	187476	8391 (1.80-1.80)
Sidechain outliers	187428	8390 (1.80-1.80)
RSRZ outliers	180081	7663 (1.80-1.80)

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 <=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	69% 27%
1	B	414	70% 26%

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
5	TRS	A	1421	-	X	-	-

2 Entry composition

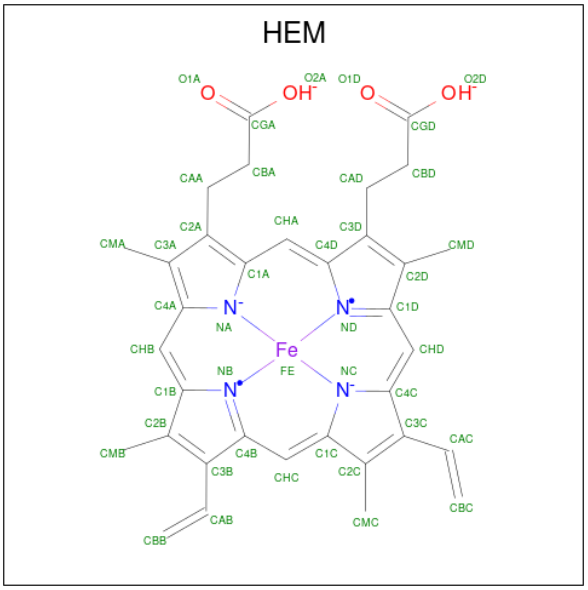
There are 7 unique types of molecules in this entry. The entry contains 7480 atoms, of which 0 are hydrogens and 0 are deuteriums.

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.

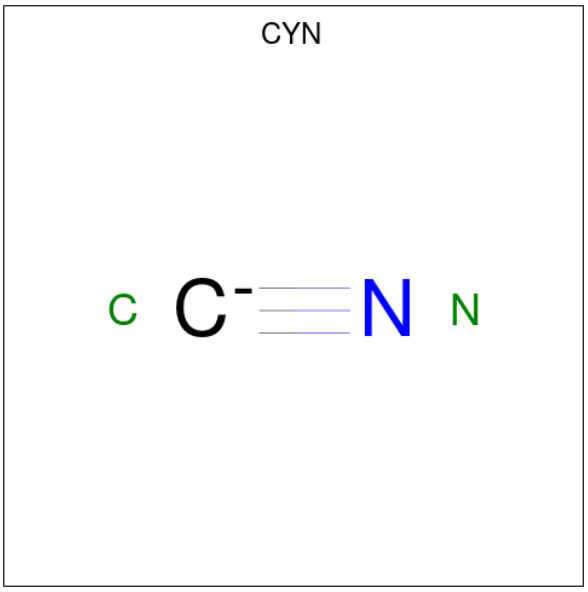
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	404	Total	C	N	O	S	0	0	0
			3199	2029	559	593	18			
1	B	405	Total	C	N	O	S	0	0	0
			3207	2033	561	595	18			

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



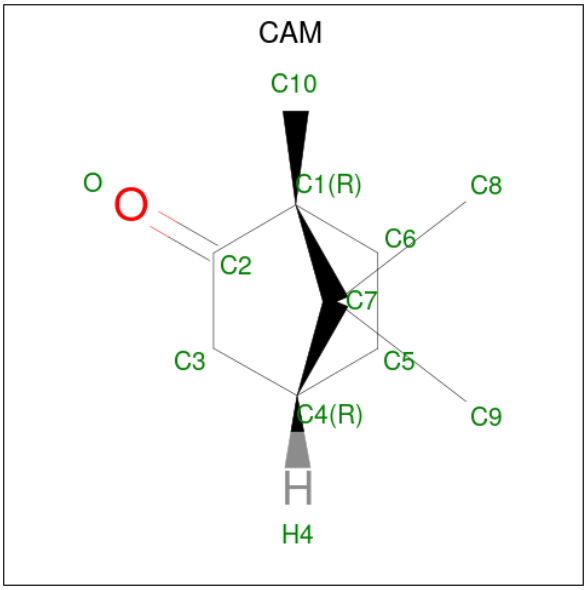
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	Fe	N	O	0	0
			43	34	1	4	4		
2	B	1	Total	C	Fe	N	O	0	0
			43	34	1	4	4		

- Molecule 3 is CYANIDE ION (CCD ID: CYN) (formula: CN).



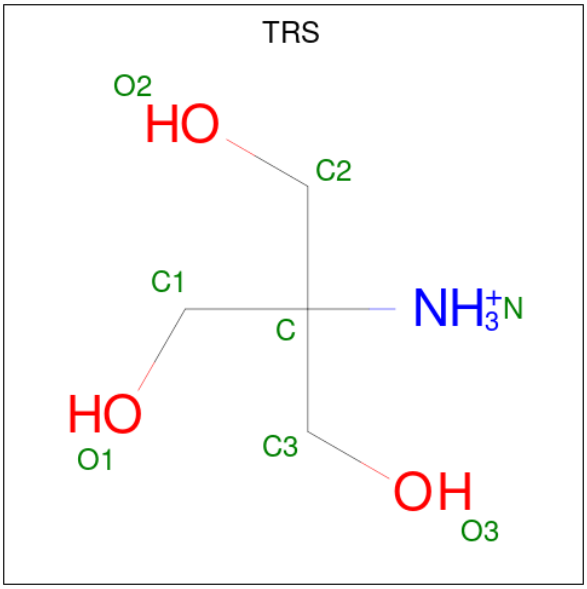
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total C N 2 1 1	0	0
3	B	1	Total C N 2 1 1	0	0

- Molecule 4 is CAMPHOR (CCD ID: CAM) (formula: $C_{10}H_{16}O$).



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

- Molecule 5 is 2-AMINO-2-HYDROXYMETHYL-PROPANE-1,3-DIOL (CCD ID: TRS) (formula: $C_4H_{12}NO_3$).



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

- Molecule 6 is POTASSIUM ION (CCD ID: K) (formula: K).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	B	3	Total	K	0	0
			3	3		

- Molecule 7 is water.

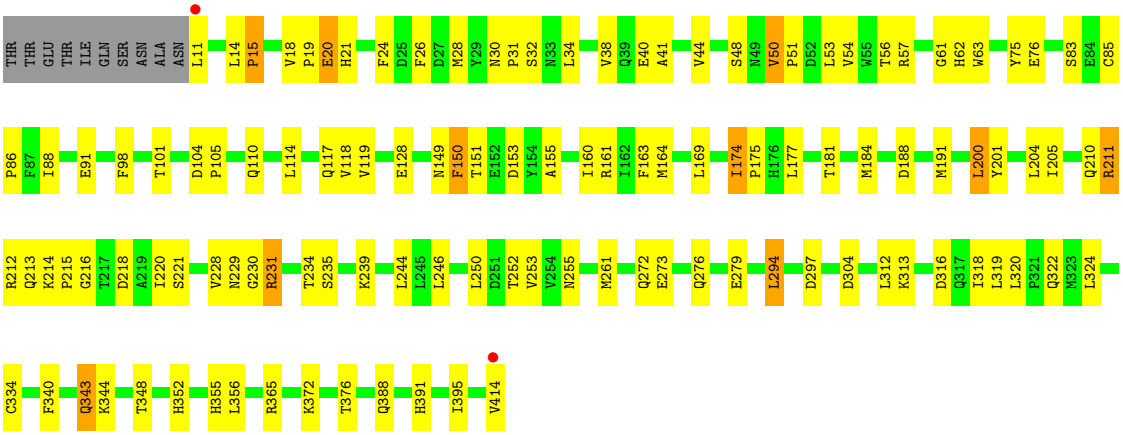
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	439	Total	O	0	0
			439	439		
7	B	512	Total	O	0	0
			512	512		

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.

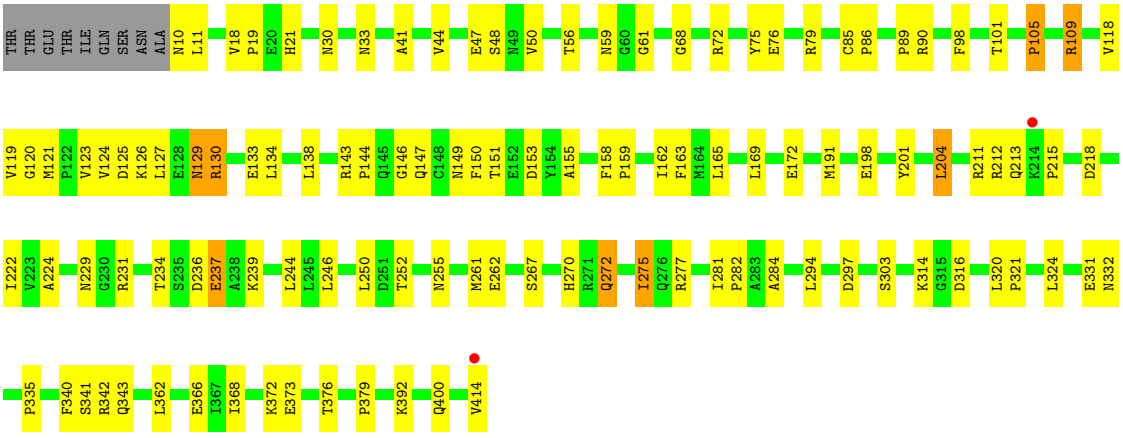
• Molecule 1: CYTOCHROME P450-CAM

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	67.00Å 61.80Å 94.60Å 90.00° 90.40° 90.00°	Depositor
Resolution (Å)	19.00 – 1.80 19.00 – 1.80	Depositor EDS
% Data completeness (in resolution range)	94.5 (19.00-1.80) 94.4 (19.00-1.80)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.08	Depositor
$< I/\sigma(I) >$ [1]	2.83 (at 1.80Å)	Xtriage
Refinement program	CNS 1.0	Depositor
R, R_{free}	0.196 , 0.236 0.193 , 0.232	Depositor DCC
R_{free} test set	3395 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å^2)	19.2	Xtriage
Anisotropy	0.414	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.37 , 46.6	EDS
L-test for twinning[2]	$< \|L\| > = 0.49$, $< L^2 > = 0.32$	Xtriage
Estimated twinning fraction	0.027 for h,-k,-l	Xtriage
F_o,F_c correlation	0.95	EDS
Total number of atoms	7480	wwPDB-VP
Average B, all atoms (Å^2)	22.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.50% of the height of the origin peak. No significant pseudotranslation is detected.*

[1] Intensities estimated from amplitudes.

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

5 Model quality (i)

5.1 Standard geometry (i)

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

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	0.79	2/3278 (0.1%)	1.18	33/4454 (0.7%)
1	B	0.72	0/3286	1.24	40/4465 (0.9%)
All	All	0.76	2/6564 (0.0%)	1.21	73/8919 (0.8%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	85	CYS	CA-C	5.85	1.58	1.52
1	A	231	ARG	CA-CB	5.01	1.60	1.54

All (73) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	B	198	GLU	CA-CB-CG	14.19	142.49	114.10
1	B	48	SER	N-CA-C	10.49	123.69	111.11
1	B	255	ASN	CA-CB-CG	-9.77	102.83	112.60
1	B	129	ASN	CA-CB-CG	-9.10	103.50	112.60
1	A	316	ASP	CA-CB-CG	8.70	121.30	112.60
1	B	105	PRO	N-CA-C	-8.66	102.16	110.47
1	B	125	ASP	CA-CB-CG	-8.62	103.98	112.60
1	A	191	MET	CA-CB-CG	8.41	130.92	114.10
1	B	79	ARG	CA-CB-CG	-8.37	97.35	114.10
1	A	211	ARG	CA-CB-CG	-8.16	97.78	114.10
1	B	324	LEU	N-CA-C	8.14	120.23	111.36
1	B	59	ASN	CA-CB-CG	-8.09	104.51	112.60
1	B	18	VAL	N-CA-C	7.88	114.73	107.56
1	A	324	LEU	N-CA-C	7.68	119.74	111.36
1	B	229	ASN	CA-CB-CG	-7.66	104.94	112.60
1	B	297	ASP	CA-CB-CG	7.66	120.26	112.60
1	B	211	ARG	CA-CB-CG	-7.57	98.97	114.10

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	B	191	MET	CA-CB-CG	7.47	129.05	114.10
1	A	297	ASP	CA-CB-CG	7.29	119.89	112.60
1	A	204	LEU	N-CA-C	7.19	120.16	111.82
1	B	252	THR	N-CA-C	6.89	118.48	110.97
1	A	20	GLU	CA-CB-CG	-6.72	100.65	114.10
1	A	150	PHE	N-CA-C	6.65	119.37	111.33
1	B	275	ILE	N-CA-C	-6.61	104.21	110.42
1	B	231	ARG	CA-CB-CG	6.60	127.29	114.10
1	B	149	ASN	CA-CB-CG	-6.53	106.07	112.60
1	B	85	CYS	CA-C-N	6.51	127.08	120.04
1	B	85	CYS	C-N-CA	6.51	127.08	120.04
1	A	188	ASP	CA-CB-CG	6.44	119.04	112.60
1	B	272	GLN	CA-CB-CG	-6.38	101.34	114.10
1	A	105	PRO	N-CA-C	-6.34	104.38	110.47
1	B	218	ASP	CA-CB-CG	-6.33	106.27	112.60
1	A	85	CYS	CA-C-N	6.32	126.86	120.04
1	A	85	CYS	C-N-CA	6.32	126.86	120.04
1	A	230	GLY	N-CA-C	6.08	120.91	113.79
1	B	109	ARG	N-CA-C	6.03	118.37	111.02
1	A	149	ASN	CA-CB-CG	-5.91	106.69	112.60
1	B	316	ASP	CA-CB-CG	5.89	118.49	112.60
1	B	130	ARG	CA-CB-CG	-5.80	102.51	114.10
1	A	250	LEU	N-CA-C	5.69	120.93	113.30
1	B	169	LEU	CA-C-N	5.69	125.97	119.83
1	B	169	LEU	C-N-CA	5.69	125.97	119.83
1	B	101	THR	N-CA-C	5.68	118.25	111.71
1	A	218	ASP	CA-CB-CG	-5.64	106.96	112.60
1	A	255	ASN	CA-CB-CG	-5.63	106.97	112.60
1	A	153	ASP	N-CA-C	5.59	119.83	112.89
1	A	252	THR	N-CA-C	5.59	117.84	111.02
1	B	262	GLU	N-CA-C	-5.58	105.10	111.07
1	A	334	CYS	CA-C-N	5.52	125.62	119.32
1	A	334	CYS	C-N-CA	5.52	125.62	119.32
1	A	50	VAL	N-CA-C	5.48	114.93	108.96
1	A	104	ASP	CA-CB-CG	5.47	118.08	112.60
1	A	304	ASP	CA-CB-CG	-5.39	107.21	112.60
1	B	343	GLN	N-CA-C	5.37	116.82	111.07
1	B	153	ASP	N-CA-C	5.36	119.82	113.28
1	A	15	PRO	CA-C-N	5.34	124.97	119.05
1	A	15	PRO	C-N-CA	5.34	124.97	119.05
1	A	174	ILE	CB-CA-C	-5.31	108.66	113.70
1	A	32	SER	N-CA-C	5.27	117.43	111.11

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	279	GLU	CA-CB-CG	5.27	124.63	114.10
1	B	237	GLU	CA-CB-CG	5.23	124.56	114.10
1	A	91	GLU	CA-CB-CG	5.21	124.51	114.10
1	A	101	THR	N-CA-C	5.20	117.69	111.71
1	B	149	ASN	N-CA-C	-5.15	98.91	108.24
1	A	48	SER	N-CA-C	5.13	117.54	111.33
1	A	229	ASN	CA-CB-CG	-5.10	107.50	112.60
1	B	204	LEU	N-CA-C	5.08	118.64	112.23
1	B	250	LEU	N-CA-C	5.08	119.58	113.17
1	B	50	VAL	CA-C-N	5.04	125.32	119.92
1	B	50	VAL	C-N-CA	5.04	125.32	119.92
1	B	47	GLU	CA-CB-CG	-5.03	104.04	114.10
1	B	109	ARG	CA-CB-CG	-5.02	104.06	114.10
1	B	198	GLU	CB-CA-C	-5.00	102.17	110.68

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	3199	0	3152	72	0
1	B	3207	0	3158	75	0
2	A	43	0	30	1	0
2	B	43	0	30	1	0
3	A	2	0	0	1	0
3	B	2	0	0	1	0
4	A	11	0	16	1	0
4	B	11	0	16	1	0
5	A	8	0	12	0	0
6	B	3	0	0	0	0
7	A	439	0	0	16	0
7	B	512	0	0	13	0
All	All	7480	0	6414	147	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 11.

All (147) 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:200:LEU:HD21	1:A:246:LEU:HD12	1.37	1.05
1:A:11:LEU:HD12	1:A:57:ARG:HB2	1.57	0.86
1:B:130:ARG:HD2	1:B:165:LEU:HD21	1.62	0.80
1:A:343:GLN:HB3	1:A:344:LYS:HD3	1.63	0.79
1:B:372:LYS:NZ	1:B:372:LYS:HB3	2.01	0.75
1:A:220:ILE:HD12	7:A:2269:HOH:O	1.87	0.73
1:B:222:ILE:HD11	7:B:2350:HOH:O	1.89	0.72
1:B:277:ARG:HH11	1:B:277:ARG:HG3	1.54	0.72
1:B:163:PHE:HE2	1:B:246:LEU:HD23	1.55	0.72
1:A:19:PRO:HB2	1:A:21:HIS:CE1	2.25	0.71
1:A:234:THR:HA	7:A:2271:HOH:O	1.90	0.71
1:B:126:LYS:HG2	1:B:130:ARG:NH2	2.07	0.68
1:B:237:GLU:CD	7:B:2373:HOH:O	2.36	0.68
1:B:163:PHE:CE2	1:B:246:LEU:HD23	2.29	0.67
1:A:212:ARG:NH1	7:A:2271:HOH:O	2.27	0.66
1:A:211:ARG:NH2	7:A:2269:HOH:O	2.30	0.65
1:B:372:LYS:HB3	1:B:372:LYS:HZ3	1.60	0.65
1:A:184:MET:HE1	1:A:200:LEU:HD12	1.81	0.63
1:B:376:THR:HG22	1:B:414:VAL:HG21	1.80	0.63
1:B:147:GLN:HG3	7:B:2271:HOH:O	1.99	0.63
2:A:1417:HEM:HMB1	2:A:1417:HEM:HBB2	1.82	0.62
1:A:62:HIS:CD2	1:A:88:ILE:HD13	2.36	0.61
1:A:128:GLU:HG3	7:A:2159:HOH:O	2.01	0.60
1:B:126:LYS:HG2	1:B:130:ARG:HH21	1.67	0.60
1:A:210:GLN:HE22	1:A:214:LYS:HD3	1.66	0.59
1:A:14:LEU:HD12	1:A:15:PRO:HD2	1.85	0.58
1:A:83:SER:O	1:A:86:PRO:HD3	2.04	0.58
1:B:234:THR:N	1:B:237:GLU:OE1	2.35	0.58
1:A:88:ILE:HD12	1:A:319:LEU:HD13	1.85	0.57
1:A:177:LEU:O	1:A:181:THR:HG23	2.05	0.57
1:B:294:LEU:H	1:B:294:LEU:HD23	1.70	0.56
1:A:201:TYR:HB3	1:A:239:LYS:HD2	1.87	0.56
1:A:114:LEU:HD12	1:A:117:GLN:NE2	2.22	0.55
1:A:98:PHE:HB3	1:A:244:LEU:HB2	1.88	0.55
1:B:400:GLN:HA	1:B:400:GLN:OE1	2.06	0.55
1:B:121:MET:HE3	1:B:124:VAL:HB	1.87	0.55
1:A:276:GLN:HE21	1:B:172:GLU:H	1.52	0.55
1:B:123:VAL:O	1:B:126:LYS:HB3	2.07	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:169:LEU:CD2	7:A:2269:HOH:O	2.56	0.54
1:B:127:LEU:HD23	1:B:130:ARG:NH2	2.22	0.54
1:A:313:LYS:HE3	7:A:2351:HOH:O	2.08	0.53
1:A:210:GLN:NE2	1:A:214:LYS:HD3	2.24	0.53
1:A:118:VAL:HG23	1:A:119:VAL:HG13	1.90	0.53
1:A:41:ALA:O	1:A:44:VAL:HG22	2.09	0.53
1:A:235:SER:N	7:A:2271:HOH:O	2.41	0.52
1:A:150:PHE:CZ	1:A:261:MET:HG3	2.45	0.52
1:A:201:TYR:O	1:A:205:ILE:HG13	2.09	0.52
1:B:368:ILE:O	1:B:372:LYS:HG3	2.10	0.52
1:B:90:ARG:NE	7:B:2176:HOH:O	2.42	0.52
1:A:169:LEU:HD21	7:A:2269:HOH:O	2.10	0.52
1:A:322:GLN:HB3	1:A:348:THR:O	2.10	0.52
1:B:11:LEU:HD23	7:B:2004:HOH:O	2.09	0.52
1:B:89:PRO:HG3	7:B:2063:HOH:O	2.09	0.52
1:B:98:PHE:HB3	1:B:244:LEU:HB2	1.91	0.51
1:A:24:PHE:HB3	1:A:54:VAL:HG21	1.92	0.50
1:A:161:ARG:NH1	7:A:2217:HOH:O	2.30	0.50
1:B:90:ARG:CZ	7:B:2176:HOH:O	2.59	0.50
3:A:1418:CYN:C	4:A:1420:CAM:H52	2.42	0.49
1:B:277:ARG:NH1	7:B:2390:HOH:O	2.45	0.49
1:B:272:GLN:NE2	1:B:272:GLN:HA	2.27	0.49
1:B:244:LEU:CD1	2:B:1417:HEM:HMD2	2.42	0.49
1:B:294:LEU:HD23	1:B:294:LEU:N	2.27	0.49
1:B:212:ARG:HG3	1:B:224:ALA:HB1	1.95	0.49
1:A:276:GLN:HE21	1:B:172:GLU:N	2.11	0.49
1:A:376:THR:CG2	1:A:414:VAL:HG21	2.42	0.49
1:B:303:SER:HA	1:B:314:LYS:HB2	1.94	0.49
1:A:34:LEU:HA	1:A:41:ALA:HB2	1.95	0.48
1:B:277:ARG:HG3	1:B:277:ARG:NH1	2.24	0.48
3:B:1418:CYN:C	4:B:1420:CAM:H52	2.43	0.48
1:A:19:PRO:HB2	1:A:21:HIS:HE1	1.76	0.48
1:B:151:THR:HA	1:B:155:ALA:HB3	1.95	0.48
1:B:158:PHE:O	1:B:162:ILE:HG12	2.13	0.48
1:A:28:MET:HE1	1:A:395:ILE:HD13	1.95	0.47
1:A:20:GLU:H	1:A:20:GLU:HG3	1.29	0.47
1:A:15:PRO:HB2	1:A:18:VAL:HG23	1.97	0.47
1:A:372:LYS:NZ	7:A:2401:HOH:O	2.46	0.47
1:A:160:ILE:O	1:A:164:MET:HG2	2.15	0.47
1:B:130:ARG:HD2	1:B:165:LEU:CD2	2.41	0.47
1:B:201:TYR:HB3	1:B:239:LYS:HD2	1.97	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:33:ASN:HB3	1:B:41:ALA:HA	1.96	0.46
1:B:72:ARG:O	1:B:76:GLU:HG3	2.15	0.46
1:B:143:ARG:HB3	1:B:144:PRO:HD3	1.98	0.46
1:B:150:PHE:CZ	1:B:261:MET:HG3	2.50	0.46
1:B:213:GLN:C	1:B:215:PRO:HD3	2.40	0.46
1:B:234:THR:HG23	1:B:237:GLU:OE1	2.16	0.46
1:B:120:GLY:O	1:B:124:VAL:HG23	2.15	0.46
1:B:281:ILE:HD12	1:B:372:LYS:HG2	1.98	0.46
1:B:158:PHE:HB3	1:B:159:PRO:CD	2.45	0.46
1:A:272:GLN:O	1:A:276:GLN:HG3	2.15	0.45
1:B:373:GLU:HA	1:B:373:GLU:OE2	2.16	0.45
1:A:163:PHE:HE2	1:A:246:LEU:CD2	2.29	0.45
1:A:151:THR:HA	1:A:155:ALA:HB3	1.99	0.45
1:A:11:LEU:CD1	1:A:57:ARG:HB2	2.39	0.45
1:A:294:LEU:HD23	1:A:294:LEU:H	1.82	0.45
1:B:143:ARG:HB3	1:B:144:PRO:CD	2.47	0.45
1:B:267:SER:OG	1:B:270:HIS:ND1	2.37	0.44
1:A:56:THR:O	1:A:61:GLY:HA2	2.17	0.44
1:B:30:ASN:ND2	7:B:2058:HOH:O	2.50	0.44
1:B:129:ASN:N	7:B:2233:HOH:O	2.50	0.44
1:B:56:THR:O	1:B:61:GLY:HA2	2.17	0.44
1:B:75:TYR:CZ	1:B:320:LEU:HB2	2.52	0.44
1:A:53:LEU:HD21	1:A:312:LEU:HD21	1.99	0.44
1:A:365:ARG:NH2	7:A:2400:HOH:O	2.50	0.44
1:B:158:PHE:HB3	1:B:159:PRO:HD3	1.99	0.44
1:A:253:VAL:HG23	7:A:2303:HOH:O	2.19	0.43
1:A:318:ILE:O	1:A:318:ILE:HG23	2.18	0.43
1:A:228:VAL:O	1:A:228:VAL:HG23	2.18	0.43
1:A:355:HIS:O	1:A:356:LEU:C	2.60	0.43
1:B:68:GLY:HA3	1:B:331:GLU:OE2	2.18	0.43
1:A:62:HIS:CG	1:A:88:ILE:HD13	2.53	0.43
1:B:392:LYS:HD2	1:B:400:GLN:OE1	2.19	0.43
1:A:344:LYS:HE2	7:A:2380:HOH:O	2.17	0.43
1:B:41:ALA:O	1:B:44:VAL:HG22	2.18	0.43
1:A:273:GLU:OE1	1:A:340:PHE:HB2	2.18	0.43
1:B:19:PRO:HB3	1:B:21:HIS:CE1	2.54	0.43
1:B:118:VAL:HG23	1:B:119:VAL:HG13	2.00	0.43
1:A:53:LEU:HD11	1:A:63:TRP:HB3	2.01	0.43
1:B:11:LEU:N	1:B:11:LEU:HD22	2.33	0.43
1:B:284:ALA:HA	1:B:340:PHE:CD1	2.54	0.42
1:A:30:ASN:ND2	7:A:2022:HOH:O	2.52	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:246:LEU:HD23	1:B:246:LEU:HA	1.85	0.42
1:B:362:LEU:O	1:B:366:GLU:HG3	2.20	0.42
1:B:341:SER:O	1:B:342:ARG:C	2.62	0.42
1:A:174:ILE:HB	1:A:175:PRO:HD3	2.02	0.42
1:A:26:PHE:CZ	1:A:31:PRO:HB3	2.55	0.42
1:A:110:GLN:H	1:A:110:GLN:HG2	1.55	0.42
1:B:163:PHE:HE2	1:B:246:LEU:CD2	2.27	0.42
1:B:204:LEU:HD21	1:B:246:LEU:HD11	2.01	0.42
1:A:276:GLN:NE2	1:B:172:GLU:H	2.17	0.41
1:B:133:GLU:HG2	7:B:2240:HOH:O	2.19	0.41
1:B:332:ASN:O	1:B:335:PRO:HD3	2.20	0.41
1:B:281:ILE:HB	1:B:282:PRO:HD3	2.01	0.41
1:A:216:GLY:H	1:A:221:SER:HB3	1.85	0.41
1:B:19:PRO:HD3	7:B:2011:HOH:O	2.19	0.41
1:A:38:VAL:HB	1:A:391:HIS:HB3	2.03	0.41
1:B:146:GLY:O	1:B:147:GLN:HB3	2.19	0.41
1:A:76:GLU:HG3	1:A:352:HIS:CD2	2.55	0.41
1:A:213:GLN:C	1:A:215:PRO:HD3	2.46	0.41
1:A:228:VAL:HG21	1:A:231:ARG:CZ	2.50	0.41
1:B:275:ILE:HD13	1:B:379:PRO:HB2	2.03	0.41
1:A:50:VAL:HA	1:A:51:PRO:HD3	1.78	0.41
1:A:163:PHE:HE2	1:A:246:LEU:HD22	1.86	0.41
1:B:105:PRO:HB3	1:B:109:ARG:CZ	2.51	0.41
1:B:236:ASP:HB3	7:B:2371:HOH:O	2.21	0.40
1:A:75:TYR:CZ	1:A:320:LEU:HB2	2.57	0.40
1:A:40:GLU:HG2	7:A:2030:HOH:O	2.22	0.40
1:A:114:LEU:O	1:A:117:GLN:HB3	2.21	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%)	390 (97%)	12 (3%)	0	100	100
1	B	403/414 (97%)	386 (96%)	16 (4%)	1 (0%)	43	31
All	All	805/828 (97%)	776 (96%)	28 (4%)	1 (0%)	48	34

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
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	349/358 (98%)	345 (99%)	4 (1%)	65	60
1	B	350/358 (98%)	346 (99%)	4 (1%)	65	60
All	All	699/716 (98%)	691 (99%)	8 (1%)	65	60

All (8) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	200	LEU
1	A	294	LEU
1	A	343	GLN
1	A	388	GLN
1	B	10	ASN
1	B	86	PRO
1	B	134	LEU
1	B	138	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (26) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	30	ASN
1	A	80	HIS

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Mol	Chain	Res	Type
1	A	110	GLN
1	A	117	GLN
1	A	132	GLN
1	A	145	GLN
1	A	210	GLN
1	A	272	GLN
1	A	276	GLN
1	A	308	HIS
1	A	317	GLN
1	A	337	HIS
1	A	343	GLN
1	A	388	GLN
1	B	21	HIS
1	B	30	ASN
1	B	33	ASN
1	B	132	GLN
1	B	147	GLN
1	B	210	GLN
1	B	213	GLN
1	B	272	GLN
1	B	276	GLN
1	B	317	GLN
1	B	337	HIS
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 10 ligands modelled in this entry, 3 are monoatomic - leaving 7 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	1417	1	50,50,50	1.26	4 (8%)	66,82,82	0.76	0
4	CAM	A	1420	-	12,12,12	0.93	1 (8%)	20,21,21	1.60	4 (20%)
4	CAM	B	1420	-	12,12,12	0.92	0	20,21,21	1.71	3 (15%)
3	CYN	B	1418	-	0,1,1	-	-	-		
3	CYN	A	1418	-	0,1,1	-	-	-		
5	TRS	A	1421	-	7,7,7	1.76	3 (42%)	9,9,9	2.05	4 (44%)
2	HEM	A	1417	1	50,50,50	1.46	7 (14%)	66,82,82	0.79	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	B	1417	1	-	3/14/54/54	-
4	CAM	A	1420	-	-	-	0/3/2/2
4	CAM	B	1420	-	-	-	0/3/2/2
5	TRS	A	1421	-	-	3/9/9/9	-
2	HEM	A	1417	1	-	2/14/54/54	-

All (15) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	1417	HEM	FE-NC	4.59	2.10	1.95
2	A	1417	HEM	CAC-C3C	-3.25	1.38	1.47
2	B	1417	HEM	CBC-CAC	3.13	1.45	1.30
2	A	1417	HEM	CBB-CAB	3.00	1.45	1.30
2	A	1417	HEM	CBC-CAC	2.95	1.44	1.30
2	B	1417	HEM	CBB-CAB	2.83	1.44	1.30
5	A	1421	TRS	O3-C3	2.83	1.51	1.42
2	B	1417	HEM	CAC-C3C	-2.78	1.39	1.47
5	A	1421	TRS	O1-C1	2.77	1.51	1.42
2	A	1417	HEM	FE-ND	2.69	2.03	1.94

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	1417	HEM	CAB-C3B	-2.56	1.40	1.47
2	B	1417	HEM	CAB-C3B	-2.56	1.40	1.47
4	A	1420	CAM	C3-C4	2.29	1.59	1.53
5	A	1421	TRS	O2-C2	2.25	1.49	1.42
2	A	1417	HEM	FE-NB	2.14	2.01	1.94

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
4	B	1420	CAM	C5-C4-C3	-5.02	92.58	106.40
4	A	1420	CAM	C5-C4-C3	-4.03	95.29	106.40
4	A	1420	CAM	C7-C1-C2	3.43	106.13	100.30
4	B	1420	CAM	C7-C1-C2	3.26	105.85	100.30
5	A	1421	TRS	C3-C-C1	3.11	120.46	110.81
5	A	1421	TRS	O1-C1-C	2.97	120.41	111.00
5	A	1421	TRS	C3-C-C2	-2.78	102.18	110.81
5	A	1421	TRS	C2-C-C1	-2.76	102.26	110.81
4	B	1420	CAM	C5-C6-C1	-2.60	99.55	104.79
4	A	1420	CAM	O-C2-C1	2.14	128.00	125.39
4	A	1420	CAM	C5-C6-C1	-2.03	100.69	104.79

There are no chirality outliers.

All (8) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	1417	HEM	C4C-C3C-CAC-CBC
5	A	1421	TRS	C1-C-C3-O3
2	A	1417	HEM	C2C-C3C-CAC-CBC
2	B	1417	HEM	C2C-C3C-CAC-CBC
2	B	1417	HEM	C4C-C3C-CAC-CBC
5	A	1421	TRS	N-C-C3-O3
5	A	1421	TRS	C2-C-C3-O3
2	B	1417	HEM	CAD-CBD-CGD-O2D

There are no ring outliers.

6 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	1417	HEM	1	0
4	A	1420	CAM	1	0
4	B	1420	CAM	1	0

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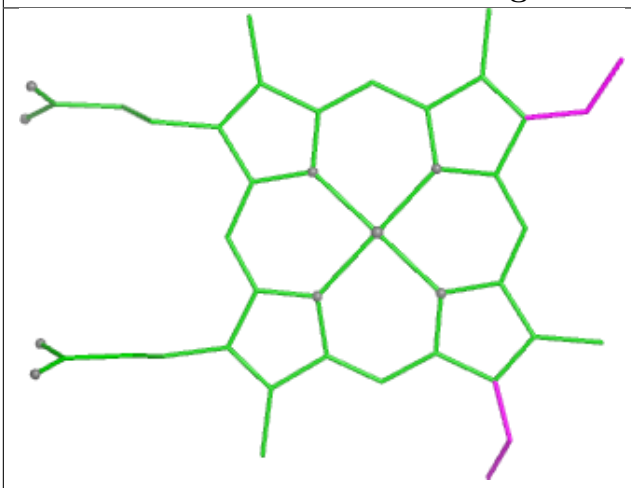
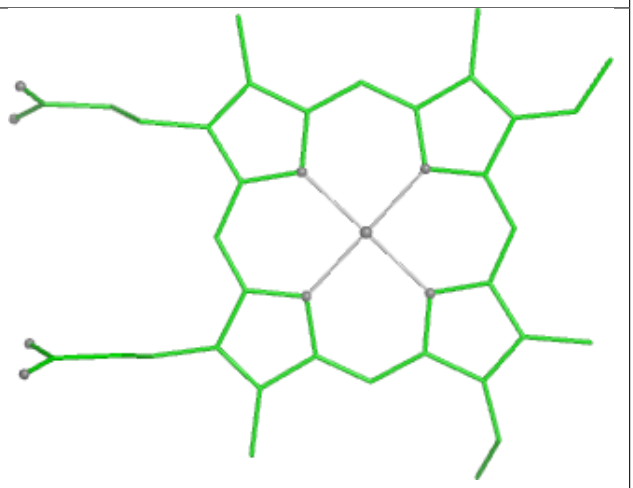
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	1418	CYN	1	0
3	A	1418	CYN	1	0
2	A	1417	HEM	1	0

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.

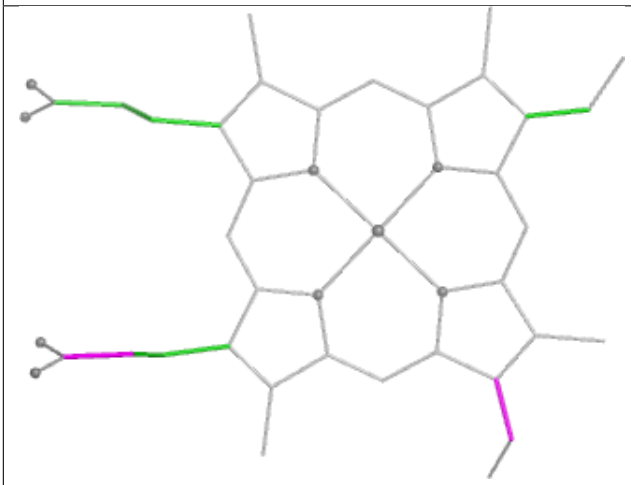
Ligand HEM B 1417



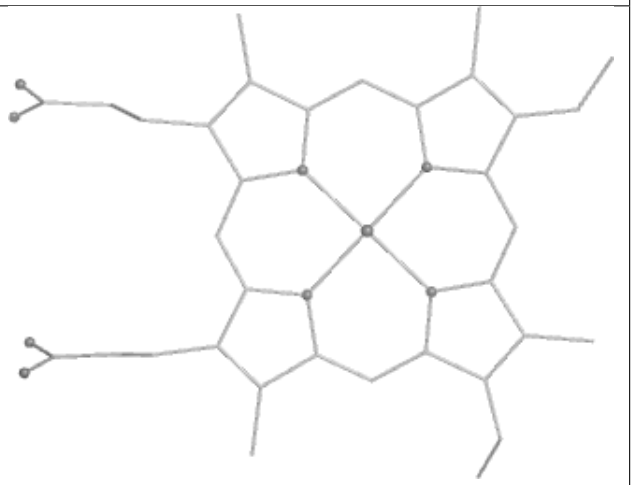
Bond lengths



Bond angles

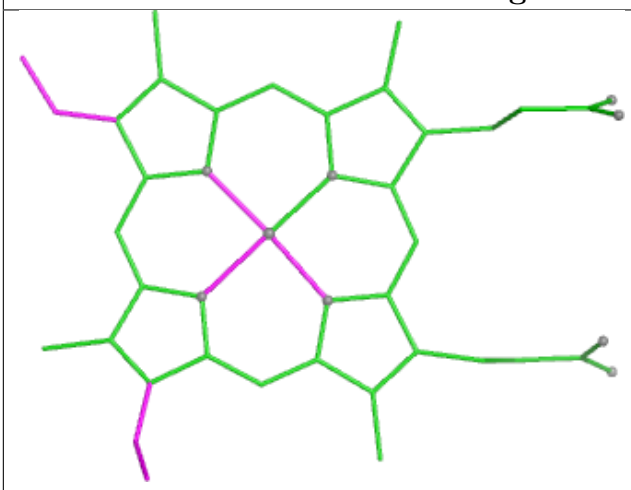


Torsions

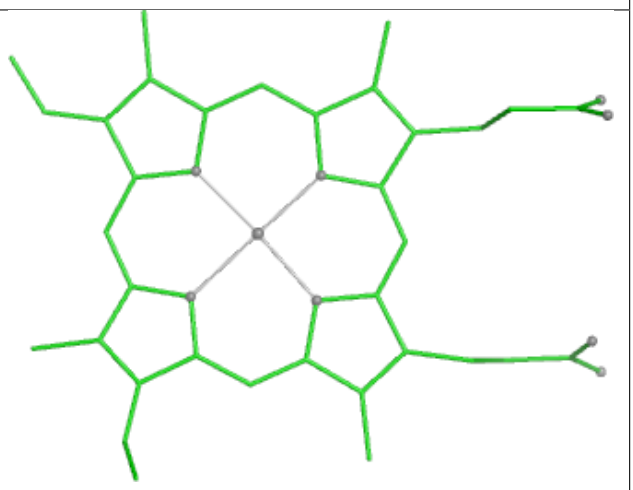


Rings

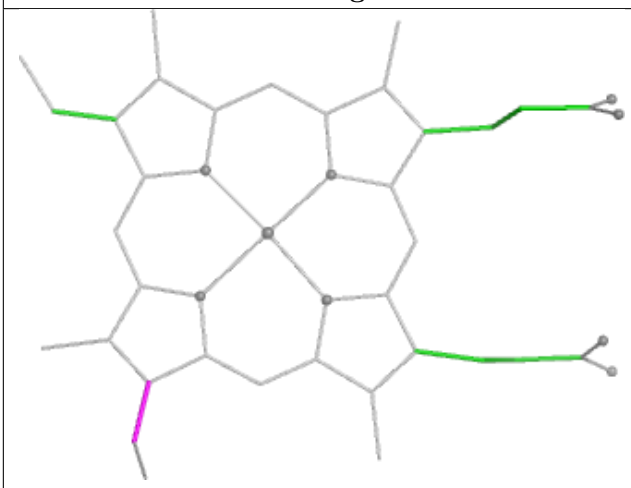
Ligand HEM A 1417



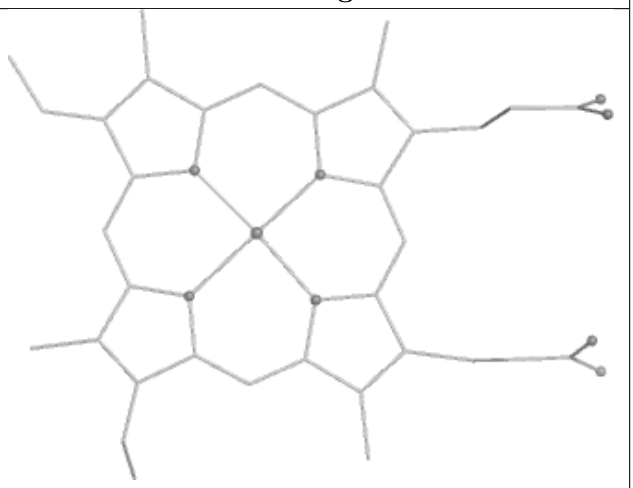
Bond lengths



Bond angles



Torsions



Rings

5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

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%)	-0.01	2 (0%)	87	88	13, 21, 36, 44	0
1	B	405/414 (97%)	-0.19	2 (0%)	87	88	10, 18, 33, 43	0
All	All	809/828 (97%)	-0.10	4 (0%)	87	88	10, 19, 35, 44	0

All (4) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	11	LEU	2.9
1	B	414	VAL	2.6
1	A	414	VAL	2.1
1	B	214	LYS	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, 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
5	TRS	A	1421	8/8	0.91	0.08	16,18,21,24	0

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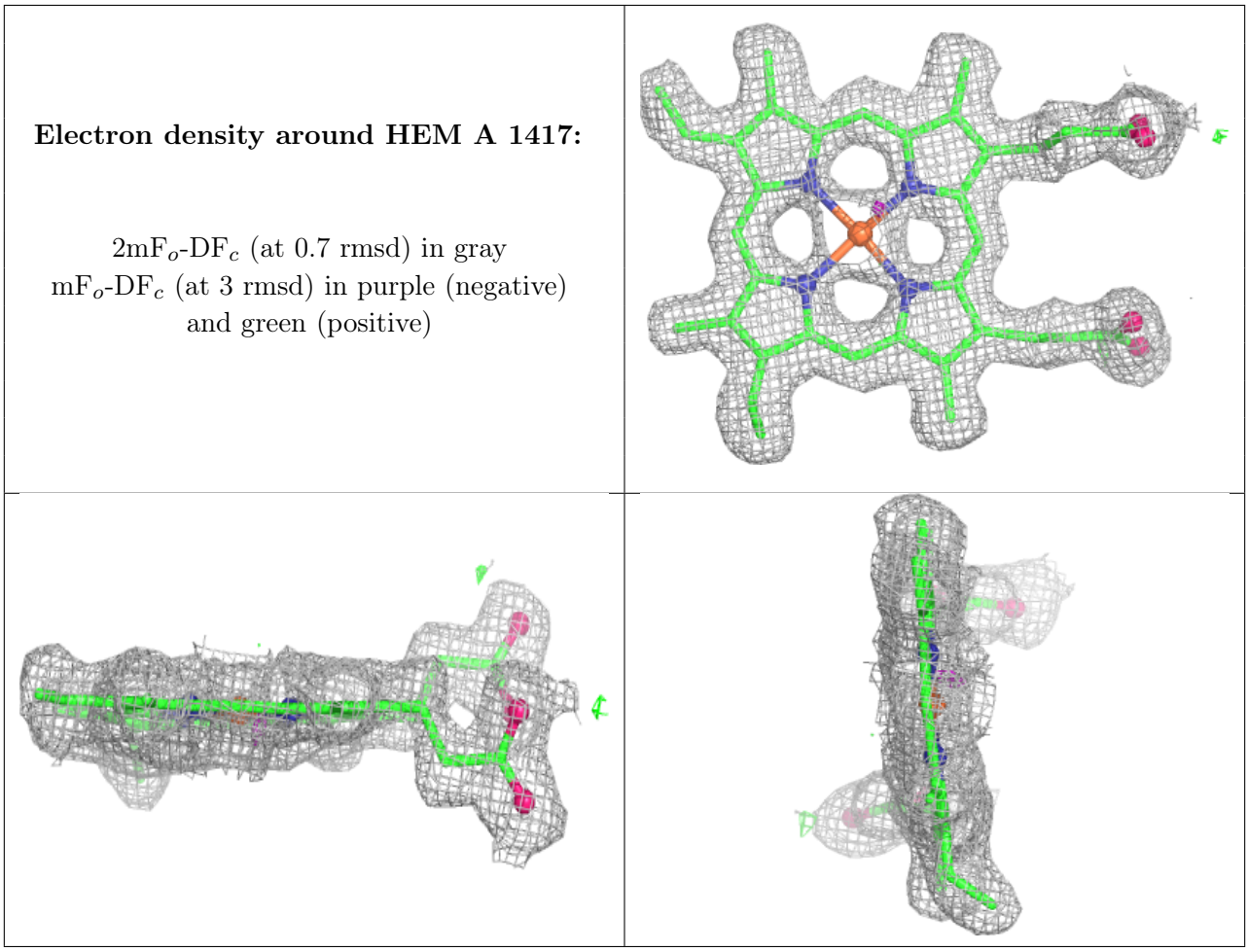
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
3	CYN	B	1418	2/2	0.92	0.14	16,16,16,17	0
4	CAM	A	1420	11/11	0.93	0.07	16,18,19,20	0
6	K	B	1419	1/1	0.95	0.04	19,19,19,19	0
3	CYN	A	1418	2/2	0.96	0.08	18,18,18,21	0
4	CAM	B	1420	11/11	0.96	0.05	10,11,13,13	0
2	HEM	A	1417	43/43	0.98	0.06	12,15,18,21	0
6	K	B	1415	1/1	0.98	0.04	11,11,11,11	0
2	HEM	B	1417	43/43	0.98	0.06	10,12,14,16	0
6	K	B	1421	1/1	0.98	0.05	18,18,18,18	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.

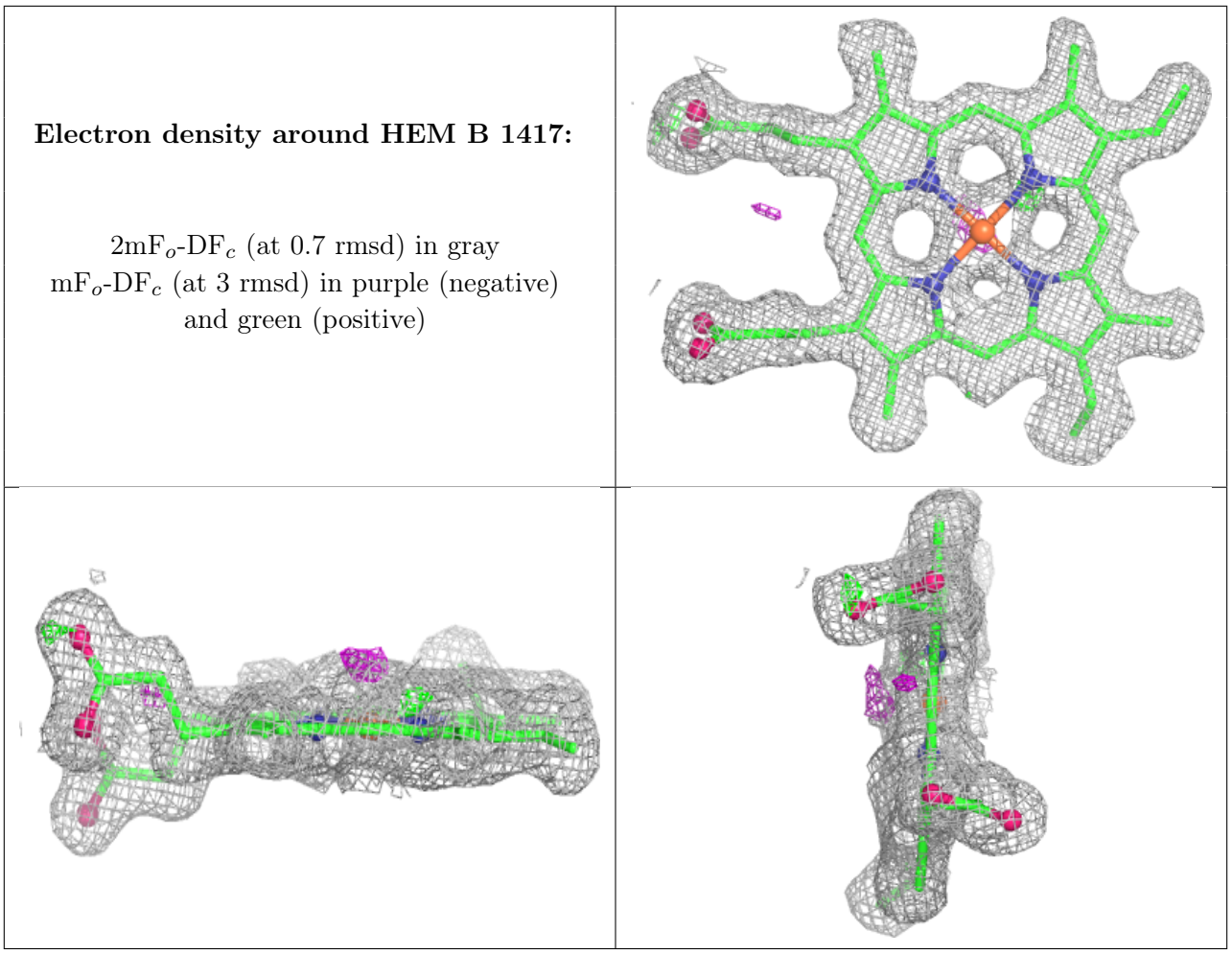
Electron density around HEM A 1417:

$2mF_o$-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 1417:

$2mF_o$-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.