



Full wwPDB X-ray Structure Validation Report ⓘ

Aug 5, 2026 – 04:03 PM JST

PDB ID : 6L39 / pdb_00006l39
Title : Cytochrome P450 107G1 (RapN)
Authors : Kim, V.C.; Kim, D.H.; Lim, Y.R.; Lee, I.H.; Lee, J.H.; Kang, L.W.
Deposited on : 2019-10-10
Resolution : 2.97 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	:	1.8.5 (274361), 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

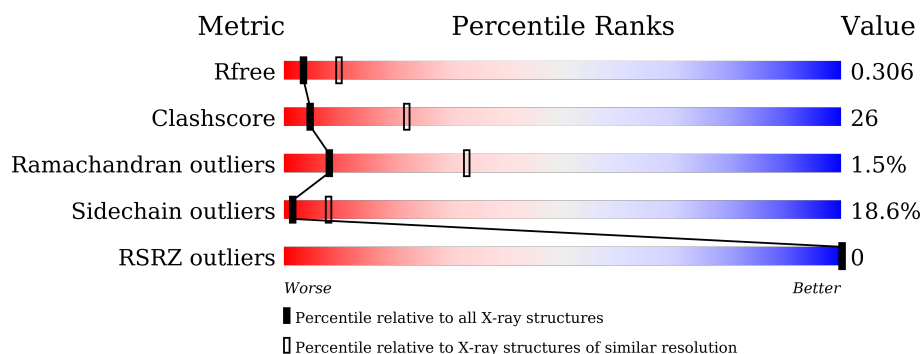
1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.97 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	3580 (3.00-2.96)
Clashscore	190562	3904 (3.00-2.96)
Ramachandran outliers	187476	3761 (3.00-2.96)
Sidechain outliers	187428	3764 (3.00-2.96)
RSRZ outliers	180081	3579 (3.00-2.96)

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	404	
1	B	404	

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	PO4	B	503	-	-	X	-

2 Entry composition [i](#)

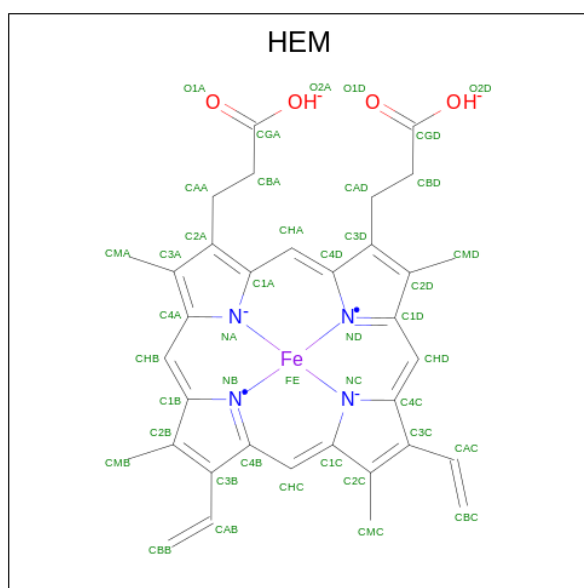
There are 5 unique types of molecules in this entry. The entry contains 6029 atoms, of which 0 are hydrogens and 0 are deuteriums.

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.

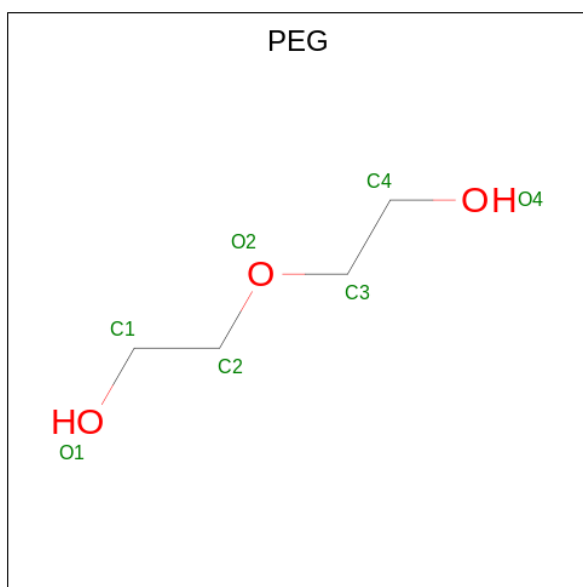
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	381	Total	C	N	O	S	0	0	0
			2965	1887	523	538	17			
1	B	378	Total	C	N	O	S	0	0	0
			2950	1875	520	540	15			

- Molecule 2 is PROTOPORPHYRIN IX CONTAINING FE (CCD ID: HEM) (formula: $C_{34}H_{32}FeN_4O_4$) (labeled as "Ligand of Interest" by depositor).



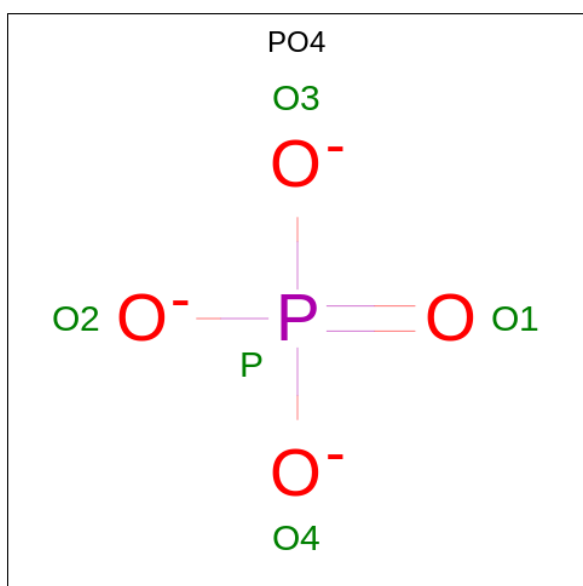
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	Fe	N	O	0	0
			43	34	1	4	4		
2	B	1	Total	C	Fe	N	O	0	0
			43	34	1	4	4		

- Molecule 3 is DI(HYDROXYETHYL)ETHER (CCD ID: PEG) (formula: $C_4H_{10}O_3$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			7	4	3		
3	A	1	Total	C	O	0	0
			7	4	3		

- Molecule 4 is PHOSPHATE ION (CCD ID: PO4) (formula: O_4P).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	B	1	Total	O	P	0	0
			5	4	1		
4	B	1	Total	O	P	0	0
			5	4	1		

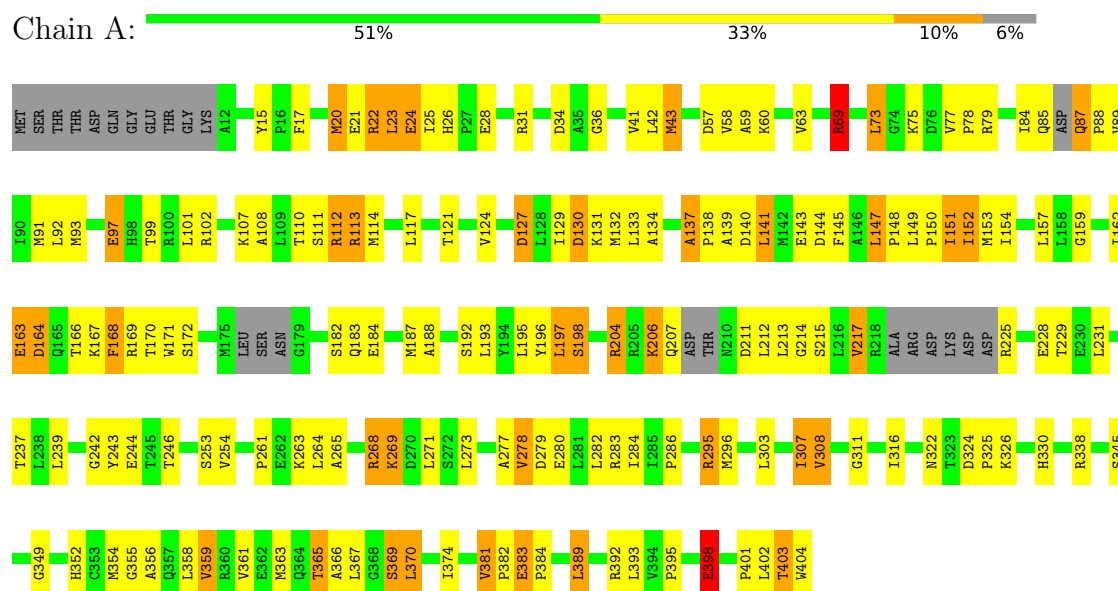
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	3	Total 3	O 3	0	0
5	B	1	Total 1	O 1	0	0

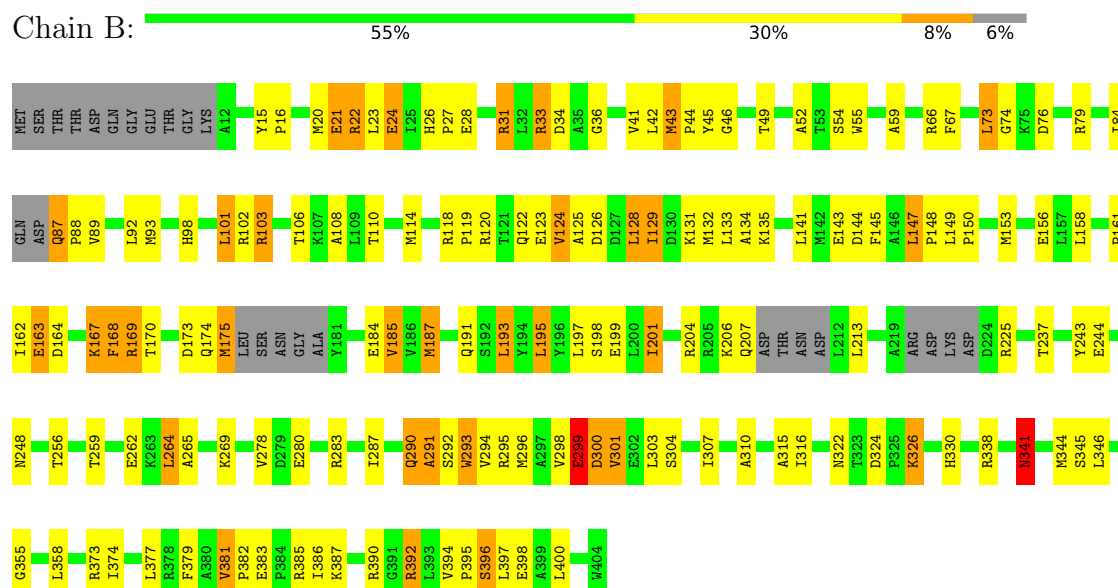
3 Residue-property plots

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



• Molecule 1: Cytochrome P450



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	59.98Å 99.80Å 62.68Å 90.00° 89.82° 90.00°	Depositor
Resolution (Å)	32.77 – 2.97 32.77 – 2.97	Depositor EDS
% Data completeness (in resolution range)	96.2 (32.77-2.97) 96.3 (32.77-2.97)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.24 (at 2.95Å)	Xtriage
Refinement program	REFMAC 5.8.0238	Depositor
R, R_{free}	0.198 , 0.313 0.209 , 0.306	Depositor DCC
R_{free} test set	997 reflections (6.51%)	wwPDB-VP
Wilson B-factor (Å ²)	72.0	Xtriage
Anisotropy	0.078	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 70.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	0.000 for -l,k,h 0.468 for h,-k,-l 0.004 for -l,-k,-h	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	6029	wwPDB-VP
Average B, all atoms (Å ²)	68.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.86% 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

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: HEM, PO4, PEG

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	0.99	0/3024	1.45	7/4105 (0.2%)
1	B	1.00	1/3009 (0.0%)	1.46	5/4084 (0.1%)
All	All	1.00	1/6033 (0.0%)	1.45	12/8189 (0.1%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	395	PRO	C-O	-5.25	1.18	1.23

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	341	ASN	CB-CA-C	9.01	119.52	110.33
1	A	273	LEU	CA-C-N	7.92	125.23	120.24
1	A	273	LEU	C-N-CA	7.92	125.23	120.24
1	B	398	GLU	N-CA-C	-7.67	103.00	111.36
1	B	125	ALA	O-C-N	5.54	128.01	122.03
1	A	198	SER	N-CA-C	-5.51	104.67	111.40
1	B	74	GLY	CA-C-N	-5.42	111.86	121.29
1	B	74	GLY	C-N-CA	-5.42	111.86	121.29
1	A	69	ARG	CA-C-N	5.18	127.53	120.54
1	A	69	ARG	C-N-CA	5.18	127.53	120.54
1	A	398	GLU	N-CA-C	-5.16	105.74	111.36
1	A	188	ALA	N-CA-C	-5.09	105.73	111.28

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	2965	0	3007	159	0
1	B	2950	0	2980	146	0
2	A	43	0	30	9	0
2	B	43	0	30	4	0
3	A	14	0	20	0	0
4	B	10	0	0	5	0
5	A	3	0	0	0	0
5	B	1	0	0	0	0
All	All	6029	0	6067	312	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 26.

All (312) 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:292:SER:CB	4:B:503:PO4:O2	1.65	1.42
1:A:163:GLU:HG2	1:A:164:ASP:OD1	1.06	1.23
1:B:292:SER:HB2	4:B:503:PO4:O2	0.99	1.16
1:A:163:GLU:CG	1:A:164:ASP:OD1	1.95	1.13
1:B:298:VAL:O	1:B:299:GLU:HB3	1.32	1.12
1:A:124:VAL:HG21	1:A:153:MET:HE3	1.31	1.08
1:B:248:ASN:ND2	1:B:396:SER:HB3	1.69	1.08
1:A:137:ALA:HB3	1:A:138:PRO:HD3	1.41	1.03
1:B:28:GLU:HA	1:B:31:ARG:HG2	1.41	1.02
1:A:148:PRO:O	1:A:152:ILE:HG13	1.61	1.00
1:A:73:LEU:HD21	1:A:92:LEU:HD13	1.43	0.99
1:A:22:ARG:HH12	1:A:383:GLU:CD	1.74	0.94
1:A:295:ARG:HG2	1:A:295:ARG:HH11	1.32	0.94
1:A:137:ALA:HB3	1:A:138:PRO:CD	1.97	0.94
1:A:112:ARG:H	1:A:112:ARG:HD2	1.33	0.93
1:A:73:LEU:HD11	1:A:88:PRO:HA	1.49	0.92
1:A:124:VAL:CG2	1:A:153:MET:HE3	1.99	0.92
1:B:248:ASN:HD21	1:B:396:SER:HB3	1.25	0.91
1:B:295:ARG:HG3	1:B:295:ARG:HH11	1.36	0.89

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:20:MET:CE	1:A:395:PRO:HB3	2.04	0.88
1:B:248:ASN:HD21	1:B:396:SER:CB	1.87	0.86
1:B:31:ARG:O	1:B:31:ARG:NH2	2.08	0.86
1:B:298:VAL:O	1:B:299:GLU:CB	2.19	0.86
1:A:22:ARG:NH1	1:A:383:GLU:OE2	2.08	0.84
1:B:292:SER:HB3	4:B:503:PO4:O2	1.76	0.84
1:A:204:ARG:HG2	1:A:204:ARG:HH21	1.42	0.82
1:B:293:TRP:HA	1:B:293:TRP:CE3	2.14	0.82
1:B:24:GLU:HA	1:B:24:GLU:OE2	1.79	0.81
1:B:26:HIS:CD2	1:B:27:PRO:HD2	2.18	0.77
1:B:89:VAL:HB	1:B:92:LEU:HD12	1.65	0.77
1:B:248:ASN:CG	1:B:396:SER:HB3	2.10	0.77
1:B:293:TRP:HA	1:B:293:TRP:HE3	1.48	0.77
1:A:307:ILE:HD12	1:A:307:ILE:N	2.02	0.75
1:B:73:LEU:HD21	1:B:92:LEU:HD13	1.69	0.75
1:A:22:ARG:NH1	1:A:383:GLU:CD	2.45	0.74
1:A:24:GLU:OE1	1:A:24:GLU:HA	1.87	0.74
1:A:73:LEU:CD2	1:A:92:LEU:HD13	2.17	0.73
1:B:31:ARG:HH21	1:B:31:ARG:HB2	1.53	0.73
1:B:295:ARG:HG3	1:B:295:ARG:NH1	1.97	0.73
1:A:271:LEU:HD12	1:A:271:LEU:N	2.03	0.73
1:A:112:ARG:H	1:A:112:ARG:CD	2.00	0.72
1:A:383:GLU:HG3	1:A:384:PRO:CD	2.20	0.72
1:B:28:GLU:CA	1:B:31:ARG:HG2	2.18	0.72
1:A:151:ILE:HG12	1:A:243:TYR:CE1	2.24	0.72
1:A:73:LEU:HD21	1:A:92:LEU:CD1	2.18	0.71
1:A:159:GLY:O	1:A:204:ARG:NH1	2.23	0.71
1:A:383:GLU:HG3	1:A:384:PRO:HD3	1.73	0.71
1:B:73:LEU:HD11	1:B:88:PRO:HA	1.72	0.71
1:B:28:GLU:O	1:B:31:ARG:N	2.24	0.70
1:A:22:ARG:NH1	1:A:383:GLU:OE1	2.24	0.70
1:A:137:ALA:CB	1:A:138:PRO:HD3	2.21	0.69
1:B:28:GLU:OE1	1:B:31:ARG:HD3	1.93	0.69
1:A:150:PRO:O	1:A:154:ILE:HG13	1.92	0.68
1:A:171:TRP:HB3	1:A:193:LEU:HD13	1.75	0.68
1:A:225:ARG:NH1	1:A:225:ARG:HG3	2.08	0.68
1:A:133:LEU:C	1:A:133:LEU:HD12	2.19	0.68
1:B:28:GLU:OE1	1:B:31:ARG:CD	2.43	0.67
1:A:295:ARG:HH11	1:A:295:ARG:CG	2.08	0.67
1:A:77:VAL:O	1:A:79:ARG:HG2	1.95	0.66
1:B:248:ASN:ND2	1:B:396:SER:CB	2.49	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:174:GLN:OE1	1:B:174:GLN:N	2.29	0.66
1:A:147:LEU:O	1:A:150:PRO:HD2	1.97	0.65
1:A:113:ARG:HH11	1:A:113:ARG:CG	2.10	0.65
1:A:113:ARG:HH11	1:A:113:ARG:HG3	1.62	0.65
1:A:363:MET:O	1:A:367:LEU:HD13	1.97	0.64
1:B:167:LYS:HB2	1:B:167:LYS:HZ1	1.62	0.64
1:B:128:LEU:N	1:B:128:LEU:HD23	2.09	0.64
1:B:28:GLU:O	1:B:31:ARG:CG	2.45	0.64
1:B:299:GLU:H	1:B:310:ALA:CB	2.11	0.63
1:B:290:GLN:H	1:B:290:GLN:NE2	1.96	0.63
1:B:76:ASP:HB2	1:B:84:ILE:HD11	1.81	0.62
1:B:167:LYS:HB2	1:B:167:LYS:NZ	2.13	0.62
1:A:113:ARG:HG3	1:A:113:ARG:NH1	2.14	0.62
1:A:102:ARG:NH1	2:A:501:HEM:O2D	2.32	0.62
1:B:76:ASP:CB	1:B:84:ILE:CD1	2.78	0.62
1:B:124:VAL:O	1:B:128:LEU:CD2	2.48	0.62
1:A:28:GLU:O	1:A:31:ARG:HB2	2.00	0.61
1:A:124:VAL:HG21	1:A:153:MET:CE	2.19	0.61
1:A:383:GLU:HG3	1:A:384:PRO:N	2.14	0.61
1:A:225:ARG:HG3	1:A:225:ARG:HH11	1.64	0.61
1:A:280:GLU:OE2	1:A:338:ARG:NE	2.32	0.61
1:A:381:VAL:HG23	1:A:382:PRO:HD2	1.82	0.61
1:A:307:ILE:HD12	1:A:307:ILE:H	1.64	0.61
2:A:501:HEM:HHA	2:A:501:HEM:HBA1	1.82	0.61
1:B:76:ASP:CB	1:B:84:ILE:HD11	2.31	0.60
1:B:287:ILE:O	1:B:394:VAL:HG23	2.01	0.60
1:A:204:ARG:HH21	1:A:204:ARG:CG	2.13	0.60
1:B:20:MET:O	1:B:20:MET:HG2	2.00	0.60
1:A:166:THR:O	1:A:167:LYS:C	2.43	0.60
1:B:26:HIS:HD2	1:B:28:GLU:H	1.50	0.59
1:B:292:SER:CA	4:B:503:PO4:O2	2.48	0.59
1:B:108:ALA:O	1:B:110:THR:HG23	2.03	0.59
1:B:392:ARG:NH1	1:B:396:SER:OG	2.35	0.59
1:A:140:ASP:HB3	1:A:143:GLU:HG2	1.85	0.59
1:A:242:GLY:HA3	2:A:501:HEM:C2C	2.38	0.59
1:B:298:VAL:O	1:B:298:VAL:HG23	2.02	0.59
1:A:23:LEU:CD2	1:A:286:PRO:HG2	2.32	0.59
1:A:108:ALA:O	1:A:113:ARG:HD2	2.02	0.59
1:B:195:LEU:N	1:B:195:LEU:HD23	2.18	0.59
1:B:169:ARG:O	1:B:170:THR:C	2.46	0.58
1:A:206:LYS:NZ	1:A:206:LYS:HB3	2.17	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:20:MET:HE1	1:A:395:PRO:HB3	1.83	0.58
1:A:164:ASP:OD1	1:A:164:ASP:N	2.33	0.58
2:A:501:HEM:HBB2	2:A:501:HEM:HHC	1.86	0.58
1:A:228:GLU:HA	1:A:231:LEU:HD12	1.86	0.57
1:A:389:LEU:HD13	1:A:398:GLU:HB2	1.85	0.57
1:B:31:ARG:HH21	1:B:31:ARG:CB	2.16	0.57
1:B:28:GLU:O	1:B:31:ARG:HG2	2.04	0.57
1:A:20:MET:O	1:A:20:MET:HG2	2.04	0.57
2:A:501:HEM:HHA	2:A:501:HEM:CBA	2.34	0.56
1:B:26:HIS:CG	1:B:27:PRO:HD2	2.41	0.56
1:B:76:ASP:HB2	1:B:84:ILE:CD1	2.36	0.56
1:B:124:VAL:O	1:B:128:LEU:HD23	2.06	0.56
1:A:366:ALA:O	1:A:370:LEU:HB2	2.05	0.56
1:A:130:ASP:O	1:A:134:ALA:HB2	2.06	0.56
1:A:355:GLY:O	1:A:358:LEU:N	2.38	0.55
1:A:69:ARG:HH22	2:A:501:HEM:CGA	2.19	0.55
1:B:300:ASP:OD1	1:B:300:ASP:N	2.29	0.55
1:B:379:PHE:CD1	1:B:386:ILE:HD11	2.42	0.55
1:A:361:VAL:O	1:A:365:THR:HG23	2.07	0.55
1:B:397:LEU:HD13	1:B:400:LEU:HB2	1.89	0.55
1:A:133:LEU:HD12	1:A:134:ALA:N	2.21	0.55
1:A:225:ARG:HH11	1:A:225:ARG:CG	2.19	0.55
1:A:367:LEU:N	1:A:367:LEU:HD12	2.22	0.55
1:B:147:LEU:N	1:B:148:PRO:HD2	2.22	0.55
1:A:295:ARG:HG2	1:A:295:ARG:NH1	2.12	0.55
1:B:193:LEU:O	1:B:197:LEU:HB2	2.06	0.54
1:A:196:TYR:CD1	1:A:196:TYR:C	2.85	0.54
1:A:73:LEU:CD1	1:A:88:PRO:HA	2.29	0.54
1:A:121:THR:HG22	1:A:153:MET:SD	2.48	0.54
1:A:57:ASP:HB3	1:A:303:LEU:HD23	1.89	0.54
2:A:501:HEM:HMC2	2:A:501:HEM:HBC2	1.89	0.54
1:A:85:GLN:NE2	1:A:87:GLN:HB2	2.23	0.53
1:B:120:ARG:O	1:B:123:GLU:HG2	2.08	0.53
1:B:15:TYR:OH	1:B:291:ALA:HB2	2.08	0.53
1:A:171:TRP:CZ2	1:A:192:SER:HB3	2.44	0.53
1:B:110:THR:O	1:B:114:MET:HG3	2.08	0.53
1:A:93:MET:CE	1:A:97:GLU:HG2	2.37	0.53
1:A:163:GLU:C	1:A:164:ASP:OD1	2.51	0.53
1:A:403:THR:OG1	1:A:404:TRP:N	2.41	0.53
1:A:147:LEU:N	1:A:148:PRO:CD	2.71	0.53
1:A:75:LYS:HB3	1:A:77:VAL:HG13	1.91	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:206:LYS:NZ	1:A:206:LYS:CB	2.73	0.52
1:B:161:PRO:O	1:B:163:GLU:N	2.42	0.52
1:B:149:LEU:HB3	1:B:150:PRO:HD3	1.91	0.52
1:B:167:LYS:NZ	1:B:167:LYS:CB	2.73	0.52
1:B:379:PHE:CE1	1:B:386:ILE:CD1	2.92	0.52
1:B:84:ILE:O	1:B:87:GLN:NE2	2.43	0.52
2:B:501:HEM:HMB1	2:B:501:HEM:HBB2	1.90	0.52
1:A:139:ALA:O	1:A:401:PRO:HA	2.09	0.52
1:B:128:LEU:N	1:B:128:LEU:CD2	2.73	0.52
1:A:168:PHE:O	1:A:171:TRP:N	2.39	0.52
1:A:330:HIS:CG	1:B:103:ARG:HD3	2.44	0.52
1:A:271:LEU:N	1:A:271:LEU:CD1	2.72	0.51
1:A:112:ARG:CD	1:A:112:ARG:N	2.72	0.51
1:A:151:ILE:HD12	1:A:169:ARG:HE	1.75	0.51
1:B:15:TYR:OH	1:B:291:ALA:N	2.43	0.51
1:A:367:LEU:N	1:A:367:LEU:CD1	2.72	0.51
1:A:193:LEU:O	1:A:197:LEU:HB2	2.11	0.51
1:A:359:VAL:HG22	2:A:501:HEM:HBB2	1.92	0.51
1:A:113:ARG:HH22	1:A:211:ASP:HA	1.76	0.50
1:B:322:ASN:ND2	1:B:345:SER:OG	2.44	0.50
1:A:28:GLU:O	1:A:31:ARG:N	2.45	0.50
1:A:296:MET:HE1	1:A:311:GLY:O	2.11	0.50
1:B:124:VAL:O	1:B:128:LEU:HG	2.11	0.50
1:B:15:TYR:CG	1:B:16:PRO:HA	2.46	0.50
1:A:23:LEU:O	1:A:23:LEU:HD22	2.11	0.50
1:A:211:ASP:O	1:A:215:SER:HB2	2.12	0.50
1:B:295:ARG:HH11	1:B:295:ARG:CG	2.15	0.50
1:A:187:MET:HE3	1:A:187:MET:HA	1.93	0.50
1:A:369:SER:O	1:A:370:LEU:C	2.55	0.50
1:A:383:GLU:CG	1:A:384:PRO:N	2.73	0.50
1:B:292:SER:HB2	4:B:503:PO4:P	2.34	0.50
1:B:355:GLY:O	1:B:358:LEU:N	2.45	0.50
1:A:308:VAL:HG22	1:A:308:VAL:O	2.12	0.50
1:A:168:PHE:O	1:A:169:ARG:C	2.55	0.49
1:B:93:MET:HE2	1:B:98:HIS:HA	1.94	0.49
1:B:299:GLU:H	1:B:310:ALA:HB2	1.76	0.49
1:A:204:ARG:HG2	1:A:204:ARG:NH2	2.18	0.49
1:A:17:PHE:HB3	1:A:26:HIS:CE1	2.46	0.49
1:A:168:PHE:O	1:A:171:TRP:HB2	2.12	0.49
1:A:295:ARG:CG	1:A:295:ARG:NH1	2.71	0.49
1:A:127:ASP:HA	1:A:130:ASP:HB2	1.95	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:114:MET:HE2	1:A:358:LEU:H	1.78	0.49
1:A:214:GLY:O	1:A:217:VAL:HG22	2.13	0.49
1:B:34:ASP:C	1:B:36:GLY:H	2.21	0.49
1:B:158:LEU:O	1:B:204:ARG:NH2	2.46	0.49
1:B:167:LYS:O	1:B:170:THR:HB	2.13	0.49
1:B:290:GLN:NE2	1:B:290:GLN:N	2.60	0.48
1:B:373:ARG:HB2	1:B:374:ILE:HD12	1.95	0.48
1:B:33:ARG:HG3	1:B:34:ASP:N	2.28	0.48
1:A:389:LEU:CD1	1:A:398:GLU:HB2	2.43	0.48
1:B:256:THR:O	1:B:259:THR:OG1	2.28	0.48
1:A:43:MET:HB3	1:A:79:ARG:O	2.14	0.48
1:B:307:ILE:N	1:B:307:ILE:HD12	2.29	0.48
1:B:341:ASN:O	1:B:341:ASN:ND2	2.46	0.48
1:B:66:ARG:CZ	1:B:301:VAL:HG13	2.44	0.48
1:B:383:GLU:C	1:B:383:GLU:CD	2.81	0.48
1:A:271:LEU:HD12	1:A:271:LEU:H	1.79	0.47
1:B:243:TYR:O	1:B:244:GLU:C	2.57	0.47
1:A:132:MET:HE2	1:A:144:ASP:HB2	1.95	0.47
1:A:133:LEU:C	1:A:133:LEU:CD1	2.88	0.47
1:A:127:ASP:OD1	1:A:127:ASP:C	2.57	0.47
1:B:93:MET:HE1	1:B:101:LEU:HD13	1.97	0.47
1:B:124:VAL:HG21	1:B:153:MET:HE3	1.96	0.47
1:A:269:LYS:HD2	1:A:269:LYS:HA	1.51	0.47
1:A:113:ARG:NH2	1:A:211:ASP:HA	2.29	0.47
1:B:129:ILE:HG22	1:B:373:ARG:NH2	2.29	0.47
1:B:141:LEU:HD11	1:B:145:PHE:CD2	2.50	0.47
1:B:330:HIS:O	1:B:338:ARG:NH2	2.47	0.47
1:A:112:ARG:HD2	1:A:112:ARG:N	2.16	0.47
1:A:34:ASP:C	1:A:36:GLY:H	2.23	0.47
1:B:41:VAL:HG23	1:B:49:THR:HG22	1.97	0.46
1:B:324:ASP:OD1	1:B:326:LYS:HG3	2.15	0.46
1:A:137:ALA:CB	1:A:138:PRO:CD	2.70	0.46
1:A:132:MET:HE3	1:A:144:ASP:HB3	1.97	0.46
1:B:45:TYR:C	1:B:46:GLY:O	2.57	0.46
1:A:307:ILE:N	1:A:307:ILE:CD1	2.73	0.46
1:B:379:PHE:CD1	1:B:386:ILE:CD1	2.99	0.46
2:B:501:HEM:HBC2	2:B:501:HEM:HHD	1.97	0.46
1:B:161:PRO:C	1:B:163:GLU:N	2.74	0.46
1:B:204:ARG:NH1	1:B:213:LEU:O	2.49	0.46
1:B:299:GLU:H	1:B:310:ALA:HB1	1.80	0.46
1:A:17:PHE:HB3	1:A:26:HIS:ND1	2.31	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:59:ALA:HB2	1:B:344:MET:HE1	1.97	0.46
1:B:248:ASN:OD1	1:B:396:SER:HB3	2.16	0.46
1:B:143:GLU:HG3	1:B:144:ASP:N	2.31	0.45
1:B:141:LEU:O	1:B:141:LEU:HG	2.13	0.45
1:A:367:LEU:CD1	1:A:367:LEU:H	2.30	0.45
1:A:355:GLY:O	1:A:356:ALA:C	2.59	0.45
1:B:184:GLU:HG3	1:B:185:VAL:HG13	1.98	0.45
1:B:293:TRP:CE3	1:B:293:TRP:CA	2.93	0.45
1:A:193:LEU:HD21	1:A:239:LEU:HD22	1.99	0.45
1:A:127:ASP:O	1:A:131:LYS:HG3	2.17	0.45
1:B:120:ARG:HG2	1:B:156:GLU:HG2	1.98	0.45
1:A:172:SER:HB3	1:A:239:LEU:HD11	1.99	0.44
1:A:283:ARG:NH2	1:A:284:ILE:HD13	2.32	0.44
1:B:122:GLN:O	1:B:126:ASP:HB2	2.18	0.44
1:A:184:GLU:O	1:A:184:GLU:HG2	2.18	0.44
1:A:324:ASP:OD1	1:A:325:PRO:HD2	2.18	0.44
1:B:153:MET:HE2	1:B:153:MET:HA	1.99	0.44
1:B:326:LYS:HB3	1:B:326:LYS:HE2	1.32	0.44
1:B:21:GLU:H	1:B:21:GLU:HG2	1.41	0.44
1:A:23:LEU:HD23	1:A:286:PRO:HG2	2.00	0.44
1:A:322:ASN:ND2	1:A:345:SER:OG	2.50	0.44
1:B:15:TYR:CD1	1:B:16:PRO:HA	2.52	0.44
1:B:381:VAL:HG22	1:B:382:PRO:CD	2.48	0.44
1:B:43:MET:HG3	1:B:44:PRO:HD2	2.00	0.44
1:B:55:TRP:CE2	1:B:344:MET:HE2	2.53	0.44
1:A:383:GLU:CD	1:A:383:GLU:C	2.85	0.44
1:A:149:LEU:HB3	1:A:150:PRO:HD3	2.00	0.43
1:A:381:VAL:CG2	1:A:382:PRO:HD2	2.48	0.43
1:A:15:TYR:HB2	1:A:41:VAL:HB	2.00	0.43
1:A:145:PHE:C	1:A:148:PRO:HD2	2.44	0.43
1:B:133:LEU:O	1:B:134:ALA:C	2.61	0.43
1:B:265:ALA:O	1:B:269:LYS:HG3	2.19	0.43
1:A:265:ALA:HA	1:A:268:ARG:HG3	1.99	0.43
1:A:141:LEU:HD13	1:A:402:LEU:HD12	2.01	0.43
1:B:264:LEU:O	1:B:264:LEU:HD23	2.18	0.43
1:B:34:ASP:C	1:B:36:GLY:N	2.77	0.43
1:B:294:VAL:HG12	1:B:315:ALA:HB2	2.00	0.43
1:A:296:MET:HE3	1:A:296:MET:HB3	1.76	0.43
1:B:164:ASP:O	1:B:168:PHE:HB3	2.19	0.43
1:B:45:TYR:HB2	1:B:79:ARG:HG3	2.01	0.43
1:B:76:ASP:HB2	1:B:79:ARG:NH1	2.34	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:117:LEU:HD11	1:A:157:LEU:HD13	2.01	0.43
1:A:204:ARG:CG	1:A:204:ARG:NH2	2.73	0.43
1:B:67:PHE:HA	1:B:296:MET:O	2.19	0.43
1:B:141:LEU:HD11	1:B:145:PHE:HD2	1.84	0.42
1:A:261:PRO:C	1:A:263:LYS:N	2.76	0.42
1:B:168:PHE:CD1	1:B:168:PHE:C	2.97	0.42
1:B:193:LEU:HD12	1:B:193:LEU:HA	1.81	0.42
1:A:187:MET:HE3	1:A:187:MET:CA	2.48	0.42
1:A:132:MET:CE	1:A:144:ASP:HB2	2.49	0.42
1:B:118:ARG:HB3	1:B:119:PRO:HD3	2.00	0.42
1:B:129:ILE:N	1:B:129:ILE:HD13	2.34	0.42
1:A:79:ARG:HG2	1:A:79:ARG:HH11	1.83	0.42
1:B:147:LEU:HA	1:B:147:LEU:HD13	1.72	0.42
1:B:201:ILE:HG13	1:B:213:LEU:HG	2.01	0.42
1:B:280:GLU:OE1	1:B:283:ARG:NE	2.45	0.42
1:A:43:MET:HB3	1:A:43:MET:HE2	1.91	0.42
1:A:381:VAL:HG23	1:A:382:PRO:CD	2.49	0.42
1:A:59:ALA:O	1:A:63:VAL:HG22	2.20	0.41
1:B:52:ALA:HB3	1:B:316:ILE:HG22	2.01	0.41
1:A:93:MET:HE3	1:A:97:GLU:HG2	2.00	0.41
1:B:73:LEU:CD1	1:B:88:PRO:HA	2.48	0.41
1:B:102:ARG:O	1:B:106:THR:HG22	2.20	0.41
1:B:22:ARG:HB2	1:B:23:LEU:H	1.68	0.41
2:B:501:HEM:HBB2	2:B:501:HEM:CMB	2.51	0.41
1:A:26:HIS:HD2	1:A:28:GLU:H	1.68	0.41
1:A:69:ARG:H	1:A:69:ARG:HG2	1.65	0.41
1:A:246:THR:HA	2:A:501:HEM:HBB1	2.03	0.41
1:B:131:LYS:O	1:B:135:LYS:HG3	2.20	0.41
1:A:277:ALA:O	1:A:278:VAL:C	2.63	0.41
1:B:381:VAL:HG22	1:B:382:PRO:HD2	2.02	0.41
1:B:26:HIS:CD2	1:B:27:PRO:CD	2.99	0.41
1:B:124:VAL:O	1:B:128:LEU:CG	2.69	0.41
1:B:93:MET:HE2	1:B:98:HIS:CA	2.51	0.41
1:B:290:GLN:HG2	1:B:291:ALA:N	2.35	0.41
1:A:58:VAL:HG22	1:A:316:ILE:HB	2.03	0.41
1:A:130:ASP:O	1:A:134:ALA:CB	2.69	0.41
2:B:501:HEM:HHA	2:B:501:HEM:HBD1	2.02	0.41
1:A:129:ILE:O	1:A:130:ASP:C	2.64	0.40
1:B:175:MET:N	1:B:175:MET:SD	2.94	0.40
1:A:63:VAL:HG12	1:A:349:GLY:HA3	2.04	0.40
1:B:187:MET:O	1:B:191:GLN:NE2	2.54	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:84:ILE:N	1:A:84:ILE:HD12	2.36	0.40
1:A:111:SER:HB2	1:A:112:ARG:HE	1.85	0.40
1:A:253:SER:O	1:A:254:VAL:C	2.64	0.40
1:B:132:MET:HG3	1:B:144:ASP:HB3	2.02	0.40
1:B:143:GLU:CG	1:B:144:ASP:N	2.85	0.40
1:B:161:PRO:C	1:B:163:GLU:H	2.29	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	371/404 (92%)	314 (85%)	50 (14%)	7 (2%)	6	27
1	B	368/404 (91%)	309 (84%)	55 (15%)	4 (1%)	11	40
All	All	739/808 (92%)	623 (84%)	105 (14%)	11 (2%)	8	33

All (11) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	162	ILE
1	B	291	ALA
1	B	299	GLU
1	A	20	MET
1	A	168	PHE
1	A	22	ARG
1	A	78	PRO
1	B	22	ARG
1	A	137	ALA
1	A	359	VAL
1	A	162	ILE

5.3.2 Protein sidechains ⓘ

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	316/343 (92%)	251 (79%)	65 (21%)	1	6
1	B	314/343 (92%)	262 (83%)	52 (17%)	2	11
All	All	630/686 (92%)	513 (81%)	117 (19%)	1	8

All (117) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	21	GLU
1	A	23	LEU
1	A	24	GLU
1	A	25	ILE
1	A	42	LEU
1	A	43	MET
1	A	60	LYS
1	A	69	ARG
1	A	73	LEU
1	A	87	GLN
1	A	89	VAL
1	A	91	MET
1	A	97	GLU
1	A	99	THR
1	A	101	LEU
1	A	107	LYS
1	A	110	THR
1	A	112	ARG
1	A	113	ARG
1	A	127	ASP
1	A	130	ASP
1	A	141	LEU
1	A	147	LEU
1	A	151	ILE
1	A	152	ILE
1	A	163	GLU
1	A	164	ASP

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Mol	Chain	Res	Type
1	A	170	THR
1	A	182	SER
1	A	183	GLN
1	A	195	LEU
1	A	197	LEU
1	A	198	SER
1	A	204	ARG
1	A	206	LYS
1	A	207	GLN
1	A	212	LEU
1	A	213	LEU
1	A	217	VAL
1	A	229	THR
1	A	237	THR
1	A	244	GLU
1	A	264	LEU
1	A	268	ARG
1	A	269	LYS
1	A	278	VAL
1	A	279	ASP
1	A	282	LEU
1	A	295	ARG
1	A	307	ILE
1	A	308	VAL
1	A	326	LYS
1	A	352	HIS
1	A	354	MET
1	A	365	THR
1	A	369	SER
1	A	370	LEU
1	A	374	ILE
1	A	381	VAL
1	A	383	GLU
1	A	389	LEU
1	A	392	ARG
1	A	393	LEU
1	A	398	GLU
1	A	403	THR
1	B	21	GLU
1	B	24	GLU
1	B	31	ARG
1	B	33	ARG

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Mol	Chain	Res	Type
1	B	42	LEU
1	B	43	MET
1	B	54	SER
1	B	73	LEU
1	B	87	GLN
1	B	101	LEU
1	B	103	ARG
1	B	124	VAL
1	B	128	LEU
1	B	129	ILE
1	B	147	LEU
1	B	163	GLU
1	B	167	LYS
1	B	168	PHE
1	B	169	ARG
1	B	173	ASP
1	B	175	MET
1	B	185	VAL
1	B	187	MET
1	B	193	LEU
1	B	195	LEU
1	B	198	SER
1	B	199	GLU
1	B	201	ILE
1	B	206	LYS
1	B	207	GLN
1	B	225	ARG
1	B	237	THR
1	B	262	GLU
1	B	264	LEU
1	B	278	VAL
1	B	290	GLN
1	B	293	TRP
1	B	299	GLU
1	B	300	ASP
1	B	301	VAL
1	B	303	LEU
1	B	304	SER
1	B	326	LYS
1	B	341	ASN
1	B	346	LEU
1	B	377	LEU

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Mol	Chain	Res	Type
1	B	381	VAL
1	B	385	ARG
1	B	387	LYS
1	B	390	ARG
1	B	392	ARG
1	B	396	SER

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	26	HIS
1	A	87	GLN
1	A	191	GLN
1	A	348	HIS
1	A	352	HIS
1	B	26	HIS
1	B	87	GLN
1	B	248	ASN
1	B	260	HIS
1	B	290	GLN
1	B	357	GLN
1	B	364	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 ⓘ

6 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
4	PO4	B	503	-	4,4,4	0.93	0	6,6,6	0.47	0
2	HEM	B	501	1	50,50,50	1.60	7 (14%)	66,82,82	2.04	23 (34%)
2	HEM	A	501	1	50,50,50	1.81	11 (22%)	66,82,82	1.92	19 (28%)
3	PEG	A	502	-	6,6,6	0.17	0	5,5,5	0.14	0
3	PEG	A	503	-	6,6,6	0.17	0	5,5,5	0.13	0
4	PO4	B	502	-	4,4,4	0.66	0	6,6,6	0.48	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	A	501	1	-	5/14/54/54	-
3	PEG	A	502	-	-	3/4/4/4	-
2	HEM	B	501	1	-	5/14/54/54	-
3	PEG	A	503	-	-	3/4/4/4	-

All (18) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	501	HEM	FE-NB	6.00	2.13	1.94
2	A	501	HEM	FE-NB	5.49	2.11	1.94
2	B	501	HEM	FE-NC	4.44	2.10	1.95
2	B	501	HEM	C1B-NB	-4.42	1.32	1.40
2	A	501	HEM	FE-NC	4.07	2.08	1.95
2	A	501	HEM	C1B-NB	-3.58	1.34	1.40
2	A	501	HEM	C3C-C4C	-3.29	1.40	1.46
2	B	501	HEM	C4D-ND	-3.11	1.34	1.40
2	A	501	HEM	C1C-C2C	-3.09	1.39	1.45
2	A	501	HEM	C4D-ND	-2.91	1.35	1.40
2	A	501	HEM	C1A-C2A	-2.74	1.39	1.44
2	B	501	HEM	C1C-C2C	-2.72	1.40	1.45

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	501	HEM	C4C-NC	-2.53	1.34	1.39
2	B	501	HEM	C4B-NB	-2.27	1.34	1.38
2	A	501	HEM	FE-ND	-2.25	1.87	1.94
2	B	501	HEM	FE-ND	-2.17	1.88	1.94
2	A	501	HEM	C3B-C4B	2.07	1.49	1.44
2	A	501	HEM	C4A-C3A	-2.02	1.38	1.43

All (42) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	501	HEM	CAA-CBA-CGA	-6.57	99.46	113.60
2	B	501	HEM	CHD-C1D-ND	5.71	130.62	124.42
2	B	501	HEM	CHC-C4B-NB	5.02	129.87	124.42
2	A	501	HEM	CHD-C1D-ND	4.25	129.03	124.42
2	B	501	HEM	CAC-C3C-C4C	4.06	134.71	124.90
2	B	501	HEM	CAD-C3D-C4D	3.81	131.31	124.66
2	A	501	HEM	C1B-NB-C4B	3.61	108.80	105.07
2	B	501	HEM	CAC-C3C-C2C	-3.44	117.27	128.60
2	B	501	HEM	C4B-C3B-C2B	-3.40	104.42	107.11
2	A	501	HEM	CHB-C1B-NB	3.25	128.38	124.37
2	B	501	HEM	CAD-C3D-C2D	-3.18	121.96	127.88
2	A	501	HEM	CHC-C4B-NB	3.16	127.85	124.42
2	A	501	HEM	CHD-C1D-C2D	-3.07	120.19	124.98
2	B	501	HEM	C4A-NA-C1A	3.07	108.35	105.35
2	A	501	HEM	C4C-NC-C1C	3.06	108.34	105.35
2	B	501	HEM	CHD-C1D-C2D	-3.00	120.29	124.98
2	A	501	HEM	CMB-C2B-C1B	2.97	129.56	125.04
2	B	501	HEM	CHA-C4D-ND	2.89	127.94	124.37
2	B	501	HEM	CAD-CBD-CGD	-2.88	107.40	113.60
2	B	501	HEM	O2D-CGD-CBD	2.74	122.83	114.03
2	B	501	HEM	CHC-C4B-C3B	-2.71	119.29	125.13
2	A	501	HEM	CAB-C3B-C2B	-2.71	119.67	128.60
2	A	501	HEM	CAA-C2A-C3A	-2.69	120.89	127.07
2	A	501	HEM	CAA-C2A-C1A	2.68	130.17	124.89
2	B	501	HEM	C4D-ND-C1D	2.66	107.82	105.07
2	A	501	HEM	C2A-C1A-NA	-2.65	107.19	110.15
2	A	501	HEM	CHB-C4A-NA	2.60	128.63	123.85
2	A	501	HEM	CHA-C4D-ND	2.55	127.52	124.37
2	A	501	HEM	C4A-NA-C1A	2.49	107.79	105.35
2	A	501	HEM	CAD-C3D-C4D	2.41	128.87	124.66
2	A	501	HEM	CHD-C4C-NC	2.35	126.98	124.44
2	B	501	HEM	O2A-CGA-CBA	2.33	121.52	114.03

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	501	HEM	C4C-NC-C1C	2.31	107.61	105.35
2	B	501	HEM	CAA-CBA-CGA	-2.27	108.71	113.60
2	A	501	HEM	CHB-C1B-C2B	-2.22	120.55	126.73
2	B	501	HEM	O1D-CGD-CBD	-2.22	115.95	123.08
2	B	501	HEM	CHB-C4A-NA	2.20	127.89	123.85
2	A	501	HEM	CAD-C3D-C2D	-2.17	123.84	127.88
2	B	501	HEM	CMC-C2C-C1C	2.07	128.35	124.71
2	B	501	HEM	CMB-C2B-C1B	2.04	128.15	125.04
2	B	501	HEM	C2A-C1A-NA	-2.04	107.87	110.15
2	B	501	HEM	CHD-C4C-C3C	2.03	128.73	125.26

There are no chirality outliers.

All (16) 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
3	A	502	PEG	O1-C1-C2-O2
2	B	501	HEM	C4D-C3D-CAD-CBD
2	B	501	HEM	C2D-C3D-CAD-CBD
2	A	501	HEM	C4B-C3B-CAB-CBB
2	B	501	HEM	C4C-C3C-CAC-CBC
3	A	503	PEG	O1-C1-C2-O2
3	A	502	PEG	C4-C3-O2-C2
3	A	503	PEG	C1-C2-O2-C3
2	A	501	HEM	CAA-CBA-CGA-O1A
3	A	503	PEG	C4-C3-O2-C2
2	A	501	HEM	CAA-CBA-CGA-O2A
2	B	501	HEM	CAA-CBA-CGA-O1A
3	A	502	PEG	O2-C3-C4-O4
2	B	501	HEM	CAA-CBA-CGA-O2A

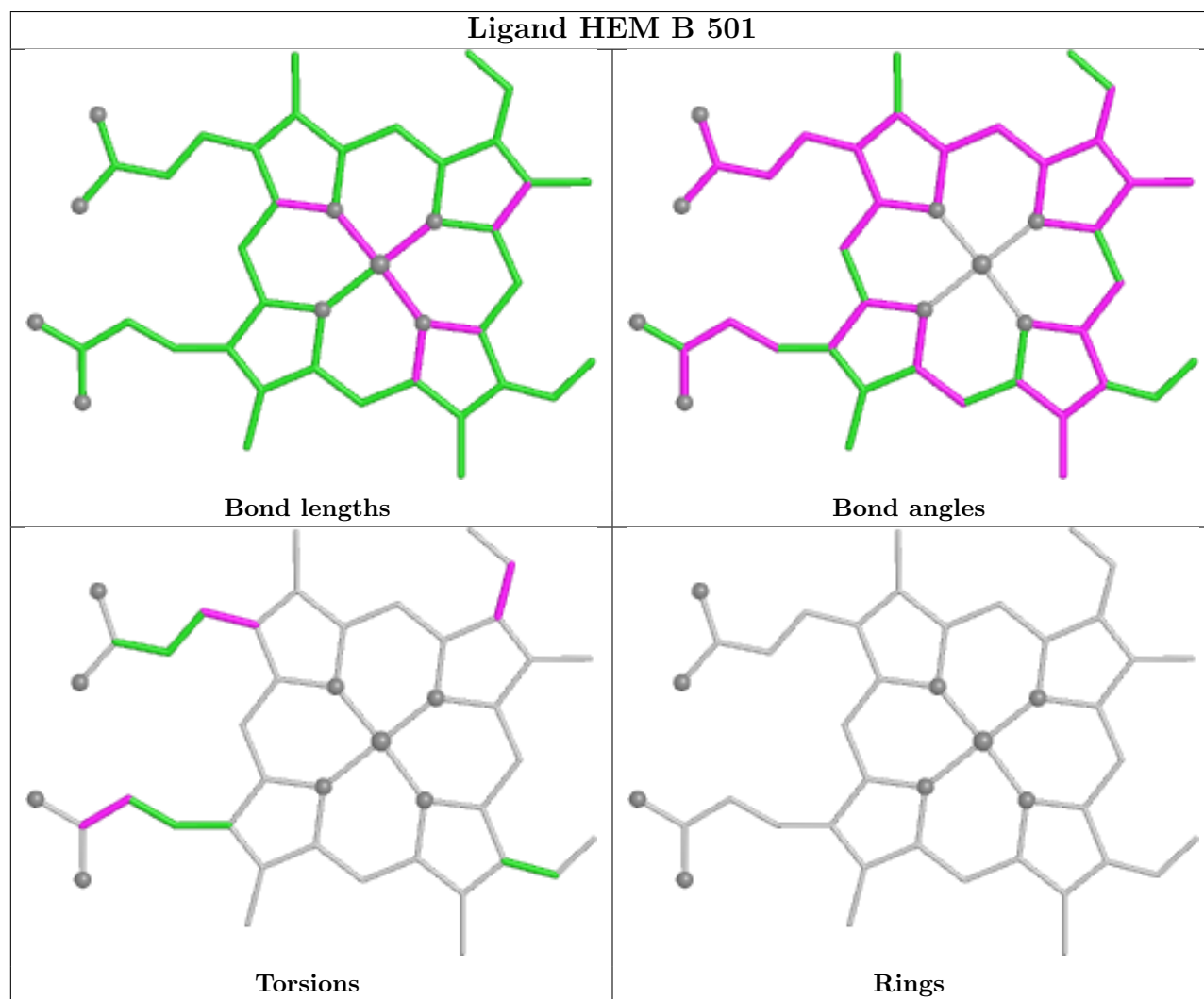
There are no ring outliers.

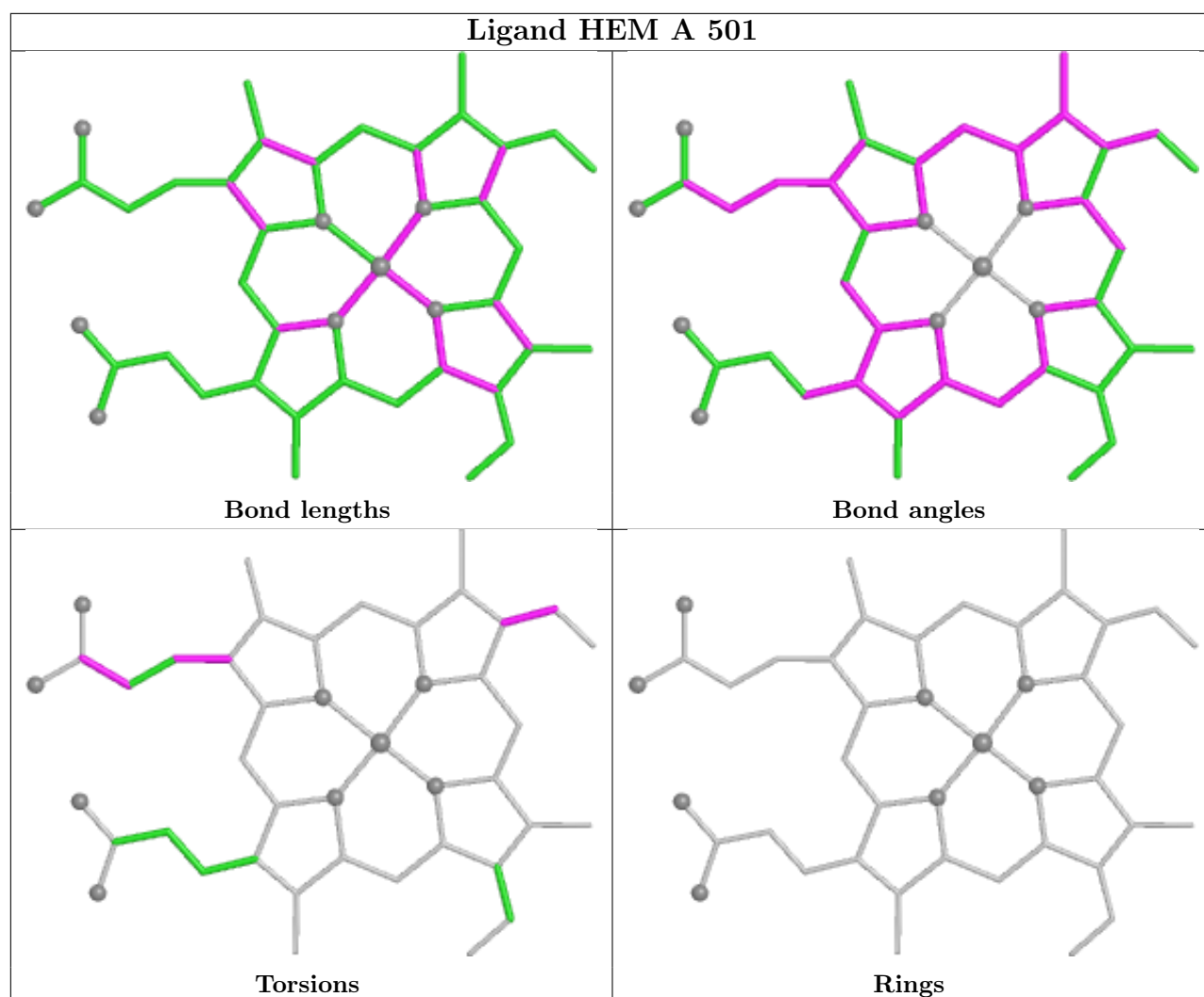
3 monomers are involved in 18 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	B	503	PO4	5	0
2	B	501	HEM	4	0
2	A	501	HEM	9	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	381/404 (94%)	-1.31	0 100 100	39, 68, 90, 105	0
1	B	378/404 (93%)	-1.28	0 100 100	37, 68, 90, 101	0
All	All	759/808 (93%)	-1.30	0 100 100	37, 68, 90, 105	0

There are no RSRZ outliers to report.

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](#)

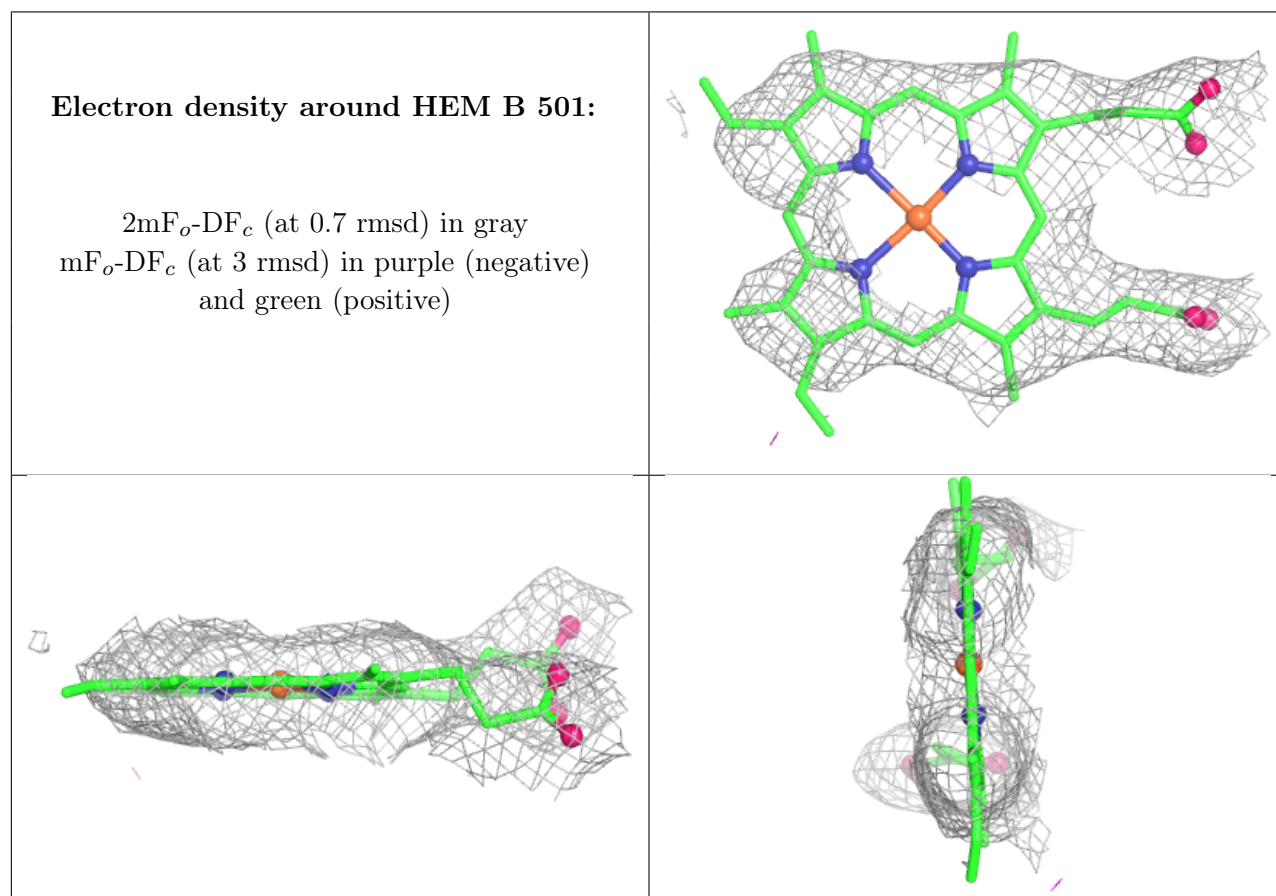
There are no oligosaccharides in this entry.

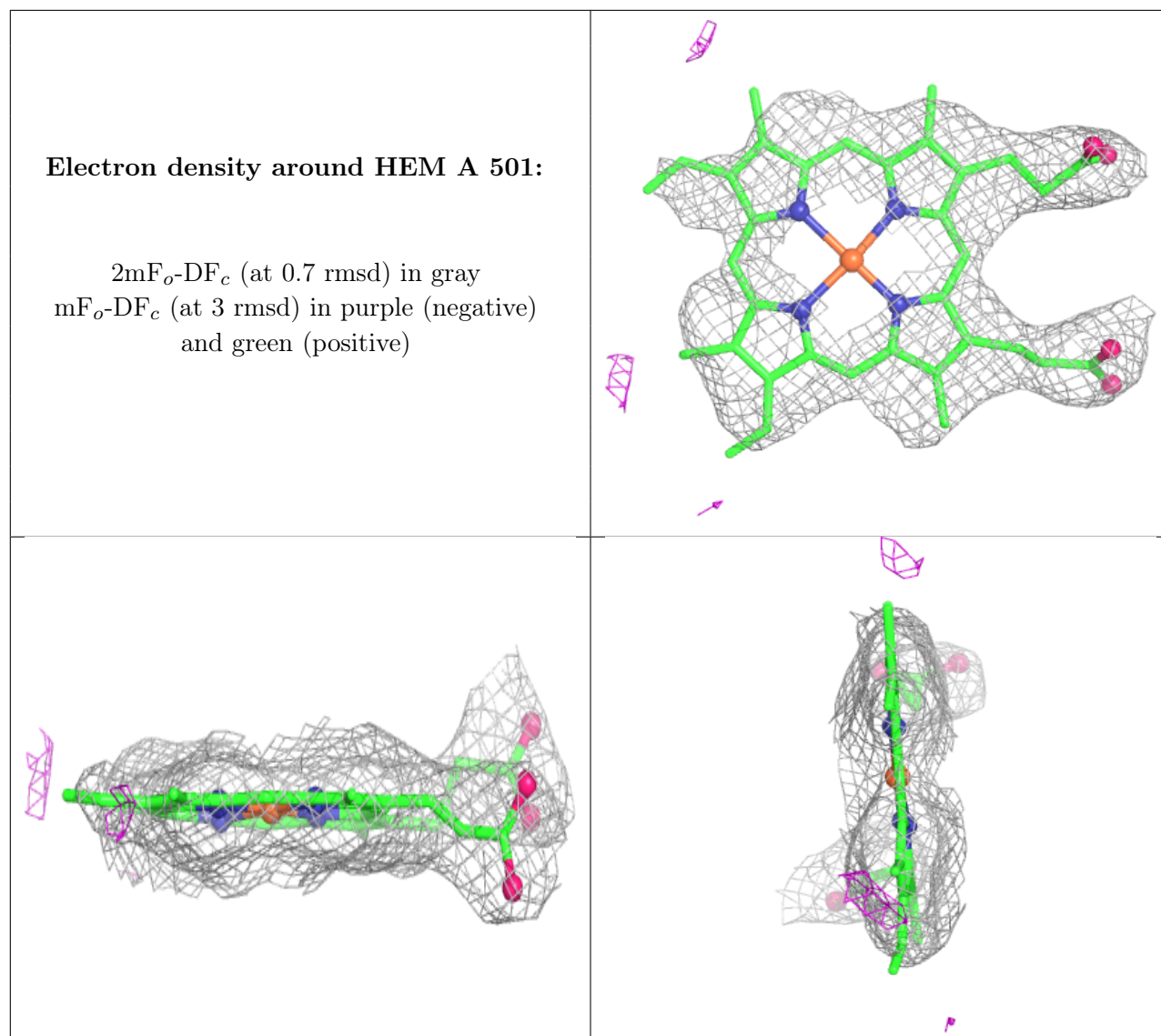
6.4 Ligands [i](#)

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
3	PEG	A	503	7/7	0.94	0.06	90,98,103,103	0
3	PEG	A	502	7/7	0.98	0.06	70,73,78,78	0
4	PO4	B	502	5/5	0.98	0.06	82,91,106,108	0
2	HEM	B	501	43/43	0.99	0.04	50,57,63,77	0
2	HEM	A	501	43/43	0.99	0.04	42,53,60,67	0
4	PO4	B	503	5/5	0.99	0.05	91,95,102,116	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.





6.5 Other polymers [i](#)

There are no such residues in this entry.