



Full wwPDB X-ray Structure Validation Report ⓘ

May 23, 2020 – 12:41 am BST

PDB ID : 1SE6
Title : Crystal Structure of Streptomyces Coelicolor A3(2) CYP158A2 from antibiotic biosynthetic pathways
Authors : Zhao, B.; Lamb, D.C.; Lei, L.; Sundaramoorthy, M.; Podust, L.M.; Waterman, M.R.
Deposited on : 2004-02-16
Resolution : 1.75 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

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

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

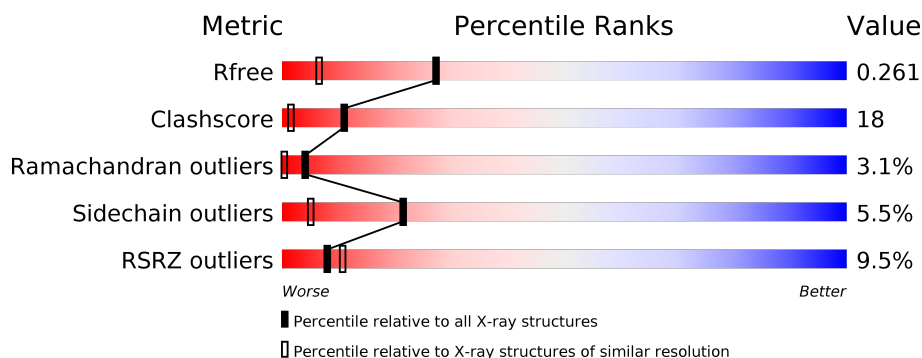
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.75 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	2340 (1.76-1.76)
Clashscore	141614	2466 (1.76-1.76)
Ramachandran outliers	138981	2437 (1.76-1.76)
Sidechain outliers	138945	2437 (1.76-1.76)
RSRZ outliers	127900	2298 (1.76-1.76)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. 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	406	<div> <div>9%</div> <div>68%</div> <div>22%</div> <div>5%</div> </div>
1	B	406	<div> <div>9%</div> <div>71%</div> <div>22%</div> <div>5%</div> </div>

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	MES	A	632	-	-	X	X

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 6734 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.

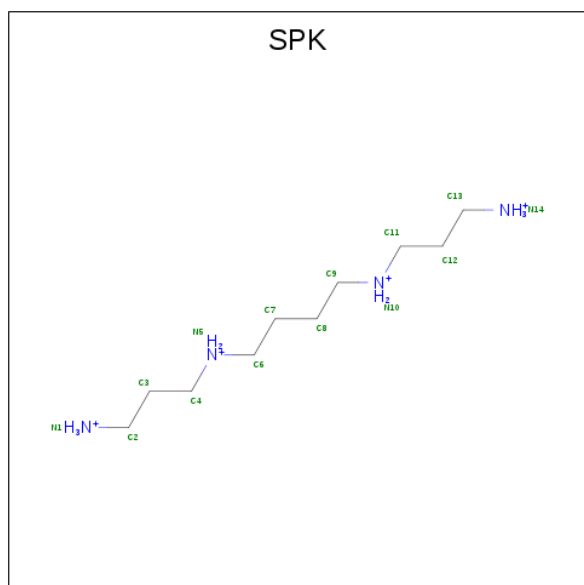
- Molecule 1 is a protein called putative cytochrome P450.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	384	Total	C	N	O	S	0	2	0
			2976	1874	543	549	10			
1	B	402	Total	C	N	O	S	0	2	0
			3107	1951	570	575	11			

There are 4 discrepancies between the modelled and reference sequences:

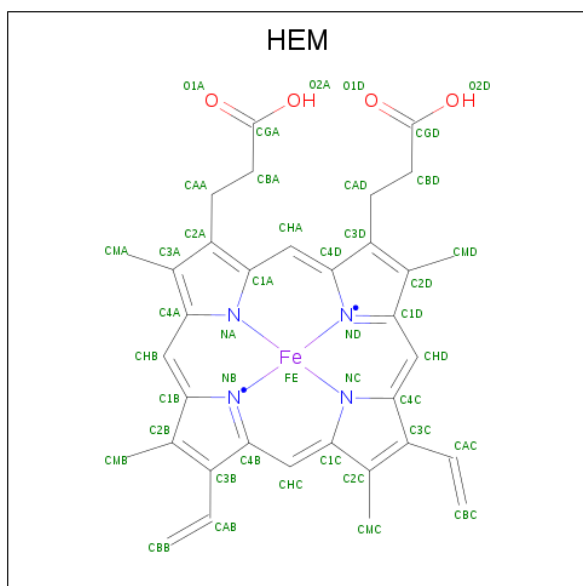
Chain	Residue	Modelled	Actual	Comment	Reference
A	405	HIS	-	CLONING ARTIFACT	UNP Q9FCA6
A	406	HIS	-	CLONING ARTIFACT	UNP Q9FCA6
B	405	HIS	-	CLONING ARTIFACT	UNP Q9FCA6
B	406	HIS	-	CLONING ARTIFACT	UNP Q9FCA6

- Molecule 2 is SPERMINE (FULLY PROTONATED FORM) (three-letter code: SPK) (formula: $C_{10}H_{30}N_4$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	N	0	0
			14	10	4		

- Molecule 3 is PROTOPORPHYRIN IX CONTAINING FE (three-letter code: HEM) (formula: $C_{34}H_{32}FeN_4O_4$).





Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	A	1	Total	C	N	O	S	0	0
			12	6	1	4	1		

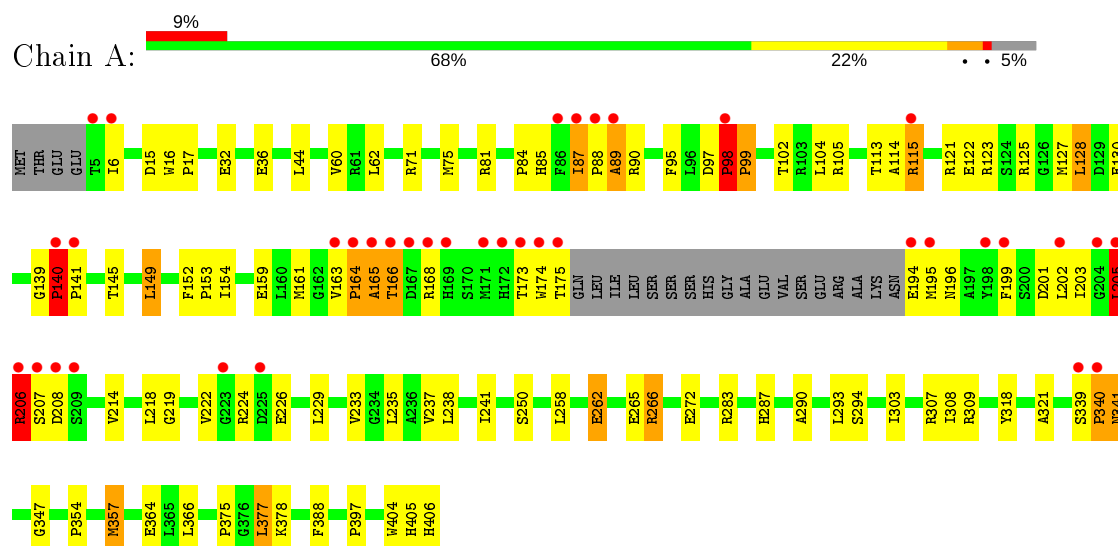
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	275	Total	O	0	0
			275	275		
5	B	264	Total	O	0	0
			264	264		

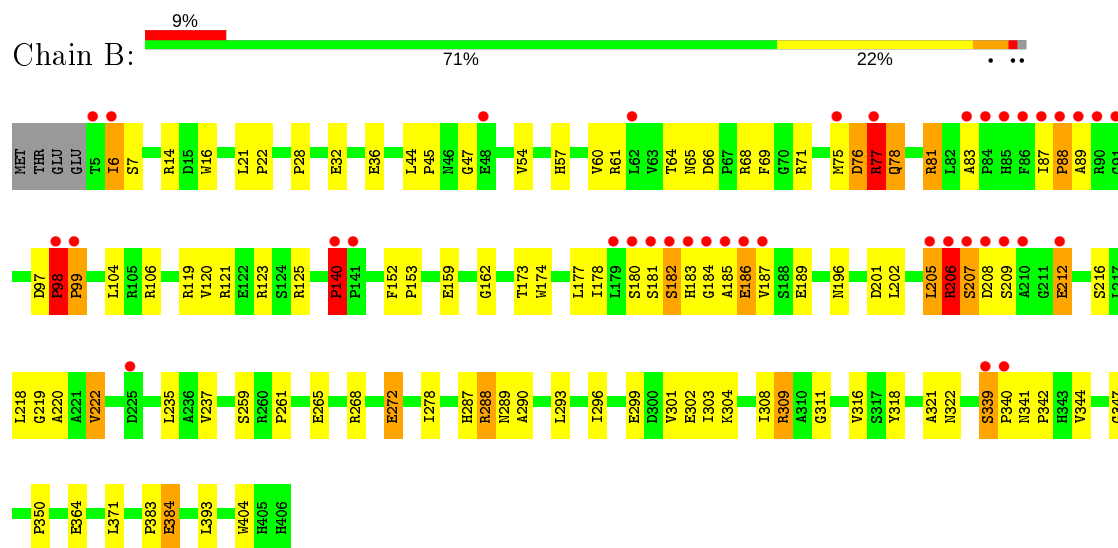
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

• Molecule 1: putative cytochrome P450



• Molecule 1: putative 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	59.55Å 79.21Å 87.43Å 90.00° 92.26° 90.00°	Depositor
Resolution (Å)	35.90 – 1.75 35.89 – 1.61	Depositor EDS
% Data completeness (in resolution range)	87.4 (35.90-1.75) 80.2 (35.89-1.61)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	0.05	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.89 (at 1.61Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.234 , 0.268 0.227 , 0.261	Depositor DCC
R_{free} test set	8715 reflections (10.03%)	wwPDB-VP
Wilson B-factor (Å ²)	11.1	Xtriage
Anisotropy	0.567	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 46.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.045 for h,-k,-l	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	6734	wwPDB-VP
Average B, all atoms (Å ²)	17.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.19% 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, SPK, MES

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	0.30	0/3045	0.73	4/4150 (0.1%)
1	B	0.35	1/3176 (0.0%)	0.68	5/4327 (0.1%)
All	All	0.33	1/6221 (0.0%)	0.70	9/8477 (0.1%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	182	SER	C-N	-6.56	1.19	1.34

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	339	SER	C-N-CD	-20.24	76.08	120.60
1	A	339	SER	C-N-CA	13.40	178.27	122.00
1	B	180	SER	N-CA-C	7.82	132.12	111.00
1	B	208	ASP	N-CA-C	-6.56	93.29	111.00
1	B	180	SER	CA-C-N	-6.53	102.84	117.20
1	A	340	PRO	CA-N-CD	-6.00	103.11	111.50
1	B	88	PRO	N-CA-CB	5.62	110.04	103.30
1	A	205	LEU	CA-CB-CG	5.60	128.19	115.30
1	B	180	SER	C-N-CA	5.29	134.92	121.70

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	2976	0	2948	112	0
1	B	3107	0	3085	111	0
2	A	14	0	30	4	0
3	A	43	0	30	0	0
3	B	43	0	30	0	0
4	A	12	0	13	6	0
5	A	275	0	0	4	0
5	B	264	0	0	10	0
All	All	6734	0	6136	223	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 18.

All (223) 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:202:LEU:CD1	1:A:206:ARG:NH2	2.01	1.23
1:A:202:LEU:HD12	1:A:206:ARG:NH2	1.56	1.18
1:B:120:VAL:HA	1:B:123:ARG:HH11	1.18	1.09
1:B:207:SER:OG	1:B:216:SER:HA	1.51	1.08
1:A:202:LEU:CD1	1:A:206:ARG:HH21	1.70	1.03
1:B:77[A]:ARG:HH11	1:B:77[A]:ARG:HB2	1.23	1.01
1:B:77[A]:ARG:HH11	1:B:77[A]:ARG:CB	1.77	0.98
1:B:185:ALA:O	1:B:189:GLU:HG3	1.65	0.96
1:B:207:SER:C	1:B:209:SER:H	1.65	0.94
1:A:206:ARG:O	1:A:206:ARG:CG	2.13	0.93
1:B:322:ASN:HD21	1:B:344:VAL:H	1.14	0.90
1:A:202:LEU:HD13	1:A:206:ARG:HH21	1.37	0.88
1:B:207:SER:OG	1:B:216:SER:CA	2.23	0.87
1:A:206:ARG:O	1:A:206:ARG:HG3	1.75	0.87
1:A:194:GLU:HG2	1:A:195:MET:H	1.39	0.85
1:A:266:ARG:HH11	1:A:266:ARG:HG3	1.45	0.81
1:B:207:SER:C	1:B:209:SER:N	2.33	0.81
1:A:201:ASP:O	1:A:205:LEU:HB3	1.81	0.81
1:A:87:ILE:HG23	1:A:88:PRO:HD3	1.60	0.81
1:B:77[B]:ARG:O	1:B:78:GLN:HB2	1.79	0.80
1:B:77[A]:ARG:O	1:B:78:GLN:HB2	1.80	0.79
1:B:212:GLU:H	1:B:212:GLU:CD	1.85	0.79
1:B:207:SER:HB2	1:B:219:GLY:HA3	1.63	0.78

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:123:ARG:HD3	1:A:159:GLU:OE2	1.84	0.78
1:A:113:THR:OG1	1:A:115:ARG:HD2	1.85	0.76
1:A:105:ARG:CZ	1:A:354:PRO:HG3	2.15	0.75
1:A:205:LEU:O	1:A:205:LEU:HD13	1.86	0.75
1:A:15:ASP:HB3	4:A:632:MES:H32	1.69	0.75
1:B:207:SER:HB3	1:B:220:ALA:N	2.03	0.73
1:B:60:VAL:HG13	5:B:608:HOH:O	1.87	0.73
1:A:194:GLU:HG2	1:A:195:MET:N	2.02	0.73
1:A:123:ARG:O	1:A:127:MET:HG3	1.89	0.72
1:A:202:LEU:HD12	1:A:206:ARG:CZ	2.19	0.72
1:A:207:SER:HB2	1:A:219:GLY:O	1.90	0.71
1:B:205:LEU:HG	1:B:206:ARG:N	2.05	0.71
1:A:115:ARG:H	1:A:115:ARG:NE	1.88	0.71
1:A:378:LYS:HZ1	2:A:631:SPK:H4A	1.56	0.71
1:B:77[A]:ARG:HH11	1:B:77[A]:ARG:CG	2.03	0.70
1:B:120:VAL:HA	1:B:123:ARG:NH1	2.01	0.70
1:B:77[A]:ARG:NH2	1:B:311:GLY:O	2.24	0.70
1:B:339:SER:CB	1:B:340:PRO:HD3	2.20	0.70
1:A:203:ILE:O	1:A:207:SER:HB3	1.92	0.69
1:B:44:LEU:HB3	1:B:45:PRO:HD2	1.75	0.68
1:B:64:THR:HG23	5:B:608:HOH:O	1.94	0.68
1:B:207:SER:O	1:B:220:ALA:HA	1.93	0.68
1:B:77[A]:ARG:HB2	1:B:77[A]:ARG:NH1	2.03	0.68
1:A:97:ASP:HB3	1:A:98:PRO:HD2	1.76	0.67
1:A:140:PRO:HB2	1:A:141:PRO:HD3	1.77	0.67
1:B:68:ARG:HE	1:B:301:VAL:HG22	1.60	0.67
1:B:339:SER:HB3	1:B:340:PRO:HD3	1.75	0.67
1:A:17:PRO:HG3	4:A:632:MES:H22	1.76	0.66
1:A:173:THR:O	1:A:175:THR:N	2.28	0.66
1:A:114:ALA:HB3	1:A:115:ARG:NH2	2.11	0.65
1:A:87:ILE:HD13	1:A:88:PRO:N	2.12	0.65
1:A:272:GLU:H	1:A:272:GLU:CD	2.00	0.65
1:A:75:MET:SD	1:A:89:ALA:O	2.55	0.65
1:A:206:ARG:O	1:A:206:ARG:HG2	1.96	0.65
1:B:66:ASP:O	1:B:68:ARG:O	2.16	0.64
1:B:119:ARG:O	1:B:123:ARG:NH1	2.31	0.63
1:B:181:SER:O	1:B:182:SER:HB3	1.99	0.62
1:B:123:ARG:HD2	1:B:159:GLU:OE2	1.99	0.62
1:A:164:PRO:HB3	5:A:886:HOH:O	1.99	0.61
1:B:104:LEU:HG	1:B:235:LEU:HD22	1.82	0.61
1:A:145:THR:HA	1:A:149:LEU:HB2	1.82	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:272:GLU:HB2	5:B:488:HOH:O	2.00	0.61
1:B:140:PRO:O	1:B:404:TRP:CZ3	2.53	0.61
1:B:81:ARG:HG3	1:B:81:ARG:HH11	1.65	0.61
1:A:195:MET:HE2	1:A:199:PHE:HE1	1.65	0.61
1:B:339:SER:CB	1:B:340:PRO:CD	2.79	0.60
1:A:196:ASN:OD1	1:A:237:VAL:HG21	2.00	0.60
1:A:121:ARG:NH2	1:A:364:GLU:OE1	2.35	0.60
1:A:87:ILE:HG23	1:A:88:PRO:CD	2.29	0.60
1:A:15:ASP:CB	4:A:632:MES:H32	2.31	0.60
1:B:184:GLY:C	1:B:186:GLU:N	2.54	0.59
1:A:115:ARG:H	1:A:115:ARG:CD	2.14	0.59
1:B:207:SER:OG	1:B:216:SER:O	2.20	0.59
1:B:218:LEU:O	1:B:222:VAL:HG13	2.03	0.58
1:B:318:TYR:CZ	1:B:347:GLY:HA2	2.38	0.58
1:B:371:LEU:HD22	5:B:667:HOH:O	2.03	0.58
1:B:121:ARG:NH1	1:B:364:GLU:OE2	2.37	0.57
1:B:69:PHE:HA	1:B:296:ILE:O	2.04	0.57
1:A:140:PRO:O	1:A:404:TRP:CZ3	2.57	0.57
1:A:378:LYS:NZ	2:A:631:SPK:H4A	2.19	0.57
1:B:162:GLY:O	1:B:206:ARG:NH2	2.29	0.56
1:A:164:PRO:O	1:A:165:ALA:HB2	2.05	0.56
1:A:152:PHE:HB3	1:A:153:PRO:HD3	1.88	0.56
1:B:207:SER:HG	1:B:216:SER:CA	2.18	0.56
1:B:316:VAL:HG11	5:B:608:HOH:O	2.06	0.55
1:A:87:ILE:HD13	1:A:87:ILE:C	2.27	0.55
1:A:194:GLU:CG	1:A:195:MET:H	2.15	0.55
1:B:68:ARG:O	1:B:69:PHE:HB2	2.05	0.55
1:A:113:THR:CB	1:A:115:ARG:HD2	2.36	0.55
1:B:206:ARG:HG3	1:B:207:SER:H	1.72	0.55
1:A:62:LEU:HD23	1:A:62:LEU:C	2.27	0.54
1:B:77[A]:ARG:NH1	1:B:77[A]:ARG:CG	2.65	0.54
1:B:152:PHE:HB3	1:B:153:PRO:HD3	1.90	0.54
1:B:207:SER:HB2	1:B:219:GLY:CA	2.35	0.54
1:A:307:ARG:NH1	1:A:309:ARG:HG3	2.23	0.54
1:A:318:TYR:CZ	1:A:347:GLY:HA2	2.43	0.54
1:B:207:SER:O	1:B:220:ALA:CA	2.56	0.54
1:A:154:ILE:HD13	1:A:168:ARG:HG3	1.90	0.54
1:B:201:ASP:O	1:B:205:LEU:HB3	2.08	0.53
1:B:99:PRO:HB2	5:B:625:HOH:O	2.06	0.53
1:A:194:GLU:OE1	1:A:194:GLU:N	2.41	0.53
1:B:212:GLU:CD	1:B:212:GLU:N	2.58	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:268:ARG:HG3	5:B:667:HOH:O	2.08	0.53
1:B:339:SER:HB2	1:B:340:PRO:HD3	1.91	0.52
1:B:384:GLU:H	1:B:384:GLU:CD	2.13	0.52
1:A:229:LEU:O	1:A:233:VAL:HG23	2.10	0.52
1:B:196:ASN:OD1	1:B:237:VAL:HG21	2.10	0.52
1:B:97:ASP:HB3	1:B:98:PRO:HD2	1.92	0.52
1:A:224:ARG:HB2	1:A:226:GLU:HG3	1.91	0.51
1:A:123:ARG:NH2	1:A:159:GLU:OE1	2.44	0.51
1:B:287:HIS:HE1	5:B:535:HOH:O	1.93	0.51
1:B:339:SER:HB2	1:B:340:PRO:CD	2.41	0.51
1:B:288:ARG:HD3	1:B:289:ASN:O	2.11	0.50
1:B:206:ARG:CG	1:B:207:SER:H	2.25	0.50
1:B:207:SER:CB	1:B:216:SER:O	2.59	0.50
1:A:87:ILE:N	1:A:88:PRO:HD2	2.25	0.50
1:A:224:ARG:NE	1:A:226:GLU:OE1	2.43	0.50
1:B:16:TRP:CZ2	1:B:290:ALA:HA	2.46	0.50
1:B:207:SER:HG	1:B:216:SER:C	2.13	0.50
1:A:114:ALA:HB3	1:A:115:ARG:CZ	2.42	0.50
1:A:168:ARG:HH11	1:A:168:ARG:HG2	1.76	0.50
1:A:75:MET:SD	1:A:90:ARG:HG2	2.52	0.50
1:A:32:GLU:O	1:A:36[B]:GLU:HG2	2.12	0.50
1:A:87:ILE:O	1:A:89:ALA:N	2.40	0.49
1:B:83:ALA:HB2	1:B:393:LEU:HD21	1.94	0.49
1:A:405:HIS:NE2	2:A:631:SPK:H3B	2.28	0.49
1:A:44:LEU:O	4:A:632:MES:H51	2.13	0.49
1:A:88:PRO:HG2	1:A:95:PHE:CE1	2.48	0.49
1:B:383:PRO:HG2	1:B:384:GLU:OE2	2.13	0.49
1:A:140:PRO:CG	1:A:406:HIS:HB2	2.43	0.49
1:A:283:ARG:HD3	1:A:341:ASN:HD21	1.78	0.49
1:B:60:VAL:HG21	1:B:321:ALA:HB2	1.95	0.49
1:A:375:PRO:O	1:A:404:TRP:HB2	2.13	0.48
1:B:212:GLU:OE2	1:B:212:GLU:N	2.40	0.48
1:B:302:GLU:O	1:B:303:ILE:HD13	2.13	0.48
1:A:98:PRO:O	1:A:102:THR:OG1	2.27	0.48
1:B:209:SER:HB2	5:B:663:HOH:O	2.14	0.48
1:A:105:ARG:NH2	1:A:354:PRO:HG3	2.28	0.47
1:A:104:LEU:HG	1:A:235:LEU:HD22	1.95	0.47
1:B:68:ARG:O	1:B:69:PHE:CB	2.63	0.47
1:A:287:HIS:HE1	5:A:798:HOH:O	1.97	0.47
1:A:140:PRO:CD	1:A:406:HIS:HB2	2.45	0.47
1:A:60:VAL:HG21	1:A:321:ALA:HB2	1.96	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:261:PRO:O	1:B:265:GLU:HG3	2.15	0.47
1:B:140:PRO:HA	1:B:404:TRP:CE2	2.49	0.47
1:B:173:THR:O	1:B:177:LEU:HG	2.15	0.47
1:B:75:MET:HE1	1:B:89:ALA:HA	1.96	0.47
1:B:207:SER:OG	1:B:216:SER:C	2.53	0.46
1:A:266:ARG:HG3	1:A:266:ARG:NH1	2.21	0.46
1:A:262:GLU:CD	1:A:262:GLU:H	2.19	0.46
1:A:266:ARG:O	1:A:266:ARG:HD3	2.16	0.46
1:A:205:LEU:O	1:A:205:LEU:HD22	2.16	0.46
1:A:114:ALA:HA	1:A:357:MET:HG2	1.97	0.46
1:A:140:PRO:HA	1:A:404:TRP:CE2	2.51	0.46
1:A:114:ALA:HA	1:A:357:MET:CG	2.46	0.46
1:A:17:PRO:HD2	5:A:761:HOH:O	2.15	0.46
1:A:161:MET:HA	1:A:214:VAL:HB	1.99	0.46
1:B:205:LEU:CG	1:B:206:ARG:N	2.71	0.46
1:A:237:VAL:O	1:A:241:ILE:HG12	2.16	0.45
1:A:262:GLU:HA	1:A:265:GLU:HG2	1.97	0.45
1:A:303:ILE:HD12	1:A:308:ILE:CD1	2.47	0.45
1:A:149:LEU:HG	1:A:250:SER:OG	2.15	0.45
1:A:165:ALA:O	1:A:166:THR:CB	2.64	0.45
1:B:318:TYR:CE2	1:B:347:GLY:HA2	2.51	0.45
1:A:71:ARG:NE	1:A:293:LEU:HD22	2.31	0.45
1:B:308:ILE:N	1:B:308:ILE:HD12	2.32	0.45
1:A:128:LEU:HD11	1:A:366:LEU:HA	1.98	0.45
1:A:233:VAL:O	1:A:237:VAL:HG23	2.17	0.45
1:A:262:GLU:HA	1:A:265:GLU:CG	2.47	0.45
1:A:16:TRP:CZ2	1:A:290:ALA:HA	2.53	0.44
1:B:81:ARG:HD3	1:B:83:ALA:O	2.17	0.44
1:B:174:TRP:O	1:B:178:ILE:HG13	2.18	0.44
1:B:341:ASN:N	1:B:342:PRO:HD3	2.31	0.44
1:B:6:ILE:HD12	1:B:7:SER:O	2.17	0.44
1:B:54:VAL:CG2	1:B:316:VAL:HG12	2.47	0.44
1:A:139:GLY:HA2	1:A:406:HIS:HD2	1.83	0.44
1:B:140:PRO:HD2	5:B:676:HOH:O	2.18	0.44
1:B:207:SER:O	1:B:219:GLY:C	2.56	0.44
1:B:202:LEU:O	1:B:206:ARG:CB	2.66	0.43
1:B:68:ARG:HE	1:B:301:VAL:CG2	2.28	0.43
1:B:181:SER:O	1:B:182:SER:CB	2.66	0.43
1:B:207:SER:CA	1:B:209:SER:H	2.29	0.43
1:B:207:SER:O	1:B:220:ALA:N	2.51	0.43
1:A:266:ARG:CG	1:A:266:ARG:NH1	2.81	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:202:LEU:HD11	1:A:206:ARG:NH2	2.18	0.43
1:A:258:LEU:HD21	1:A:377:LEU:HD13	2.01	0.43
1:A:17:PRO:HG3	4:A:632:MES:C2	2.45	0.43
1:A:218:LEU:O	1:A:222:VAL:HG23	2.19	0.43
1:A:115:ARG:NH1	5:A:700:HOH:O	2.52	0.42
1:A:97:ASP:C	1:A:99:PRO:HD2	2.39	0.42
1:A:15:ASP:CG	4:A:632:MES:H32	2.39	0.42
1:B:71:ARG:CZ	1:B:293:LEU:HD23	2.50	0.42
1:B:106:ARG:HG3	1:B:106:ARG:HH11	1.84	0.42
1:B:57:HIS:CE1	1:B:61:ARG:HD3	2.54	0.42
1:A:388:PHE:HA	1:A:397:PRO:HA	2.01	0.42
1:A:84:PRO:O	1:A:85:HIS:ND1	2.53	0.42
1:B:21:LEU:HA	1:B:22:PRO:HD3	1.84	0.42
1:B:309:ARG:HD3	1:B:309:ARG:N	2.34	0.42
1:A:114:ALA:CA	1:A:357:MET:HG2	2.49	0.42
1:A:309:ARG:HD3	1:A:309:ARG:H	1.83	0.42
1:A:378:LYS:HZ1	2:A:631:SPK:H7B	1.84	0.42
1:B:259:SER:O	1:B:261:PRO:HD3	2.20	0.42
1:A:113:THR:HB	1:A:115:ARG:HH11	1.85	0.41
1:A:115:ARG:NE	1:A:115:ARG:N	2.63	0.41
1:B:120:VAL:CA	1:B:123:ARG:HH11	2.08	0.41
1:B:183:HIS:CB	1:B:187:VAL:CG1	2.99	0.41
1:B:28:PRO:O	1:B:32:GLU:HG3	2.20	0.41
1:B:278:ILE:CD1	1:B:364:GLU:HA	2.50	0.41
1:B:207:SER:HB3	1:B:220:ALA:H	1.82	0.41
1:B:76:ASP:O	1:B:77[A]:ARG:CB	2.68	0.41
1:A:88:PRO:O	1:A:89:ALA:O	2.38	0.41
1:B:207:SER:HB3	1:B:216:SER:O	2.21	0.41
1:B:65:ASN:HD22	1:B:350:PRO:HD3	1.86	0.41
1:B:68:ARG:HD2	1:B:299:GLU:CG	2.51	0.40
1:B:47:GLY:CA	1:B:81:ARG:NH1	2.84	0.40
1:A:293:LEU:HD23	1:A:294:SER:N	2.37	0.40
1:B:184:GLY:O	1:B:185:ALA:C	2.60	0.40
1:A:173:THR:C	1:A:175:THR:N	2.75	0.40
1:A:81:ARG:NH2	1:A:85:HIS:HA	2.37	0.40
1:B:68:ARG:NE	1:B:301:VAL:HG22	2.32	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	382/406 (94%)	359 (94%)	11 (3%)	12 (3%)	4	0
1	B	402/406 (99%)	369 (92%)	20 (5%)	13 (3%)	4	0
All	All	784/812 (97%)	728 (93%)	31 (4%)	25 (3%)	4	0

All (25) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	89	ALA
1	A	164	PRO
1	A	166	THR
1	A	206	ARG
1	A	208	ASP
1	A	340	PRO
1	B	78	GLN
1	B	87	ILE
1	B	88	PRO
1	B	186	GLU
1	B	207	SER
1	B	339	SER
1	A	98	PRO
1	B	77[A]	ARG
1	B	77[B]	ARG
1	B	99	PRO
1	B	140	PRO
1	A	6	ILE
1	A	99	PRO
1	A	140	PRO
1	A	174	TRP
1	B	98	PRO
1	B	205	LEU
1	B	206	ARG
1	A	165	ALA

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	315/336 (94%)	296 (94%)	19 (6%)	19	4
1	B	329/336 (98%)	311 (94%)	18 (6%)	21	5
All	All	644/672 (96%)	607 (94%)	37 (6%)	21	5

All (37) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	87	ILE
1	A	98	PRO
1	A	115	ARG
1	A	122	GLU
1	A	125	ARG
1	A	128	LEU
1	A	130[A]	GLU
1	A	130[B]	GLU
1	A	140	PRO
1	A	149	LEU
1	A	163	VAL
1	A	205	LEU
1	A	206	ARG
1	A	238	LEU
1	A	262	GLU
1	A	266	ARG
1	A	341	ASN
1	A	357	MET
1	A	377	LEU
1	B	6	ILE
1	B	14	ARG
1	B	36	GLU
1	B	76	ASP
1	B	77[A]	ARG
1	B	77[B]	ARG
1	B	81	ARG
1	B	98	PRO

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Mol	Chain	Res	Type
1	B	125	ARG
1	B	140	PRO
1	B	206	ARG
1	B	212	GLU
1	B	222	VAL
1	B	272	GLU
1	B	288	ARG
1	B	304	LYS
1	B	309	ARG
1	B	384	GLU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (11) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	65	ASN
1	A	240	GLN
1	A	249	ASN
1	A	287	HIS
1	A	341	ASN
1	A	406	HIS
1	B	65	ASN
1	B	240	GLN
1	B	252	GLN
1	B	287	HIS
1	B	322	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry

4 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	MES	A	632	-	12,12,12	0.45	0	14,16,16	0.75	0
3	HEM	A	430	1,5	27,50,50	2.07	8 (29%)	17,82,82	1.79	5 (29%)
3	HEM	B	430	1,5	27,50,50	2.11	6 (22%)	17,82,82	1.87	6 (35%)
2	SPK	A	631	-	13,13,13	0.53	0	12,12,12	1.19	1 (8%)

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	MES	A	632	-	-	3/6/14/14	0/1/1/1
3	HEM	A	430	1,5	-	0/6/54/54	-
3	HEM	B	430	1,5	-	0/6/54/54	-
2	SPK	A	631	-	-	5/11/11/11	-

All (14) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	430	HEM	C3B-CAB	4.18	1.56	1.47
3	A	430	HEM	C3C-C2C	-4.06	1.34	1.40
3	B	430	HEM	C3C-C2C	-4.04	1.34	1.40
3	A	430	HEM	C3B-CAB	3.96	1.56	1.47
3	A	430	HEM	CBB-CAB	3.84	1.54	1.29
3	A	430	HEM	CBC-CAC	3.80	1.54	1.29
3	B	430	HEM	CBC-CAC	3.76	1.54	1.29
3	B	430	HEM	CBB-CAB	3.75	1.54	1.29
3	B	430	HEM	C3C-CAC	3.70	1.55	1.47
3	A	430	HEM	C3C-CAC	3.62	1.55	1.47
3	B	430	HEM	CAA-C2A	2.78	1.56	1.52

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	430	HEM	CAA-C2A	2.10	1.55	1.52
3	A	430	HEM	C1C-C2C	2.04	1.47	1.42
3	A	430	HEM	C3B-C2B	-2.04	1.37	1.40

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	430	HEM	CBA-CAA-C2A	3.69	119.30	112.49
3	A	430	HEM	CBA-CAA-C2A	3.50	118.93	112.49
2	A	631	SPK	C3-C4-N5	2.92	120.01	112.14
3	A	430	HEM	CMA-C3A-C4A	-2.88	124.04	128.46
3	B	430	HEM	CMA-C3A-C4A	-2.81	124.15	128.46
3	B	430	HEM	CMB-C2B-C3B	2.62	129.58	124.68
3	A	430	HEM	CMD-C2D-C1D	-2.58	124.50	128.46
3	B	430	HEM	CMC-C2C-C3C	2.50	129.35	124.68
3	B	430	HEM	CMD-C2D-C1D	-2.50	124.62	128.46
3	A	430	HEM	CMB-C2B-C3B	2.44	129.24	124.68
3	A	430	HEM	CMC-C2C-C3C	2.43	129.23	124.68
3	B	430	HEM	CBD-CAD-C3D	-2.16	108.50	112.48

There are no chirality outliers.

All (8) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	632	MES	C8-C7-N4-C3
2	A	631	SPK	C2-C3-C4-N5
2	A	631	SPK	N1-C2-C3-C4
2	A	631	SPK	C11-C12-C13-N14
4	A	632	MES	C8-C7-N4-C5
2	A	631	SPK	C6-C7-C8-C9
4	A	632	MES	C7-C8-S-O3S
2	A	631	SPK	C7-C6-N5-C4

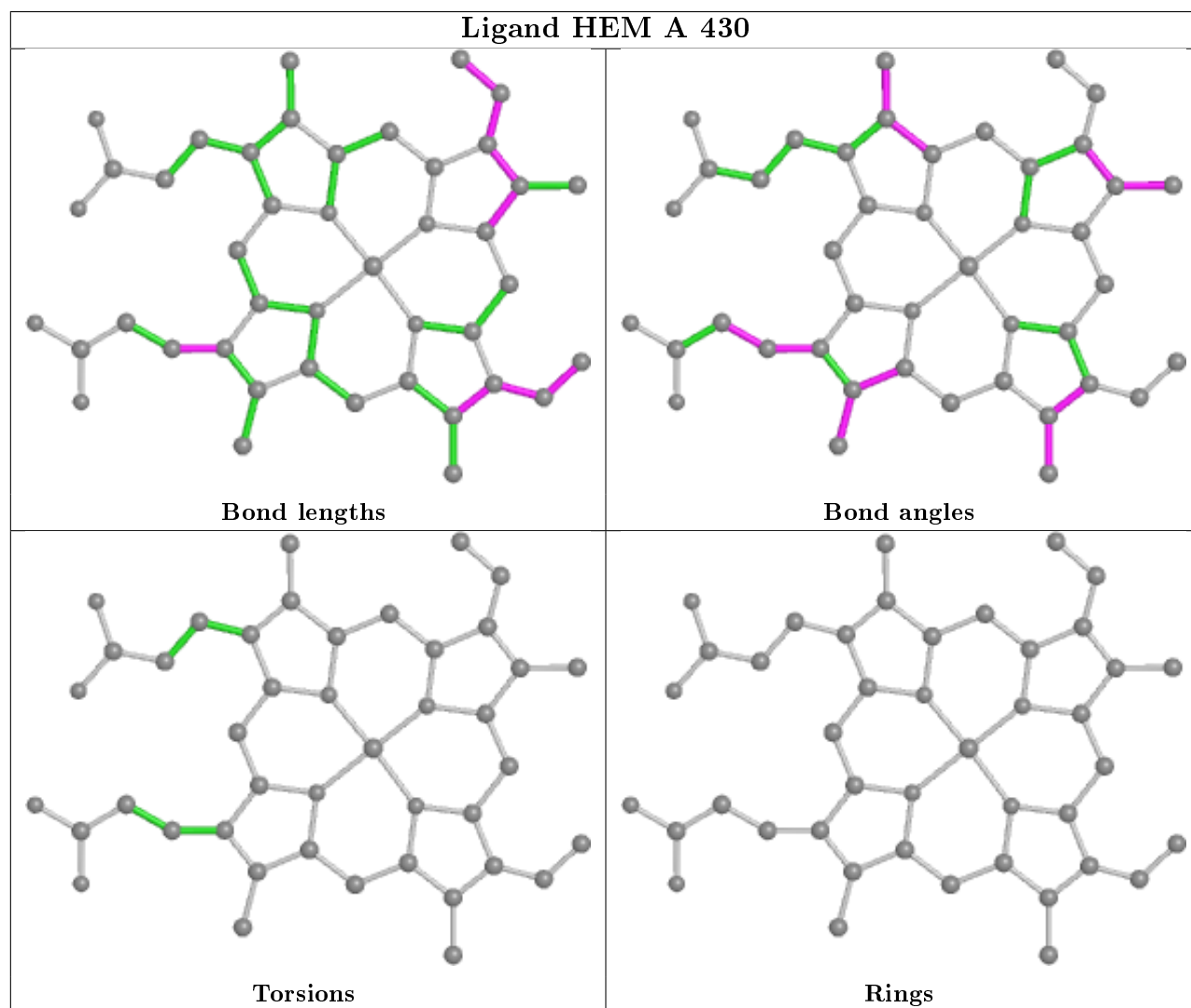
There are no ring outliers.

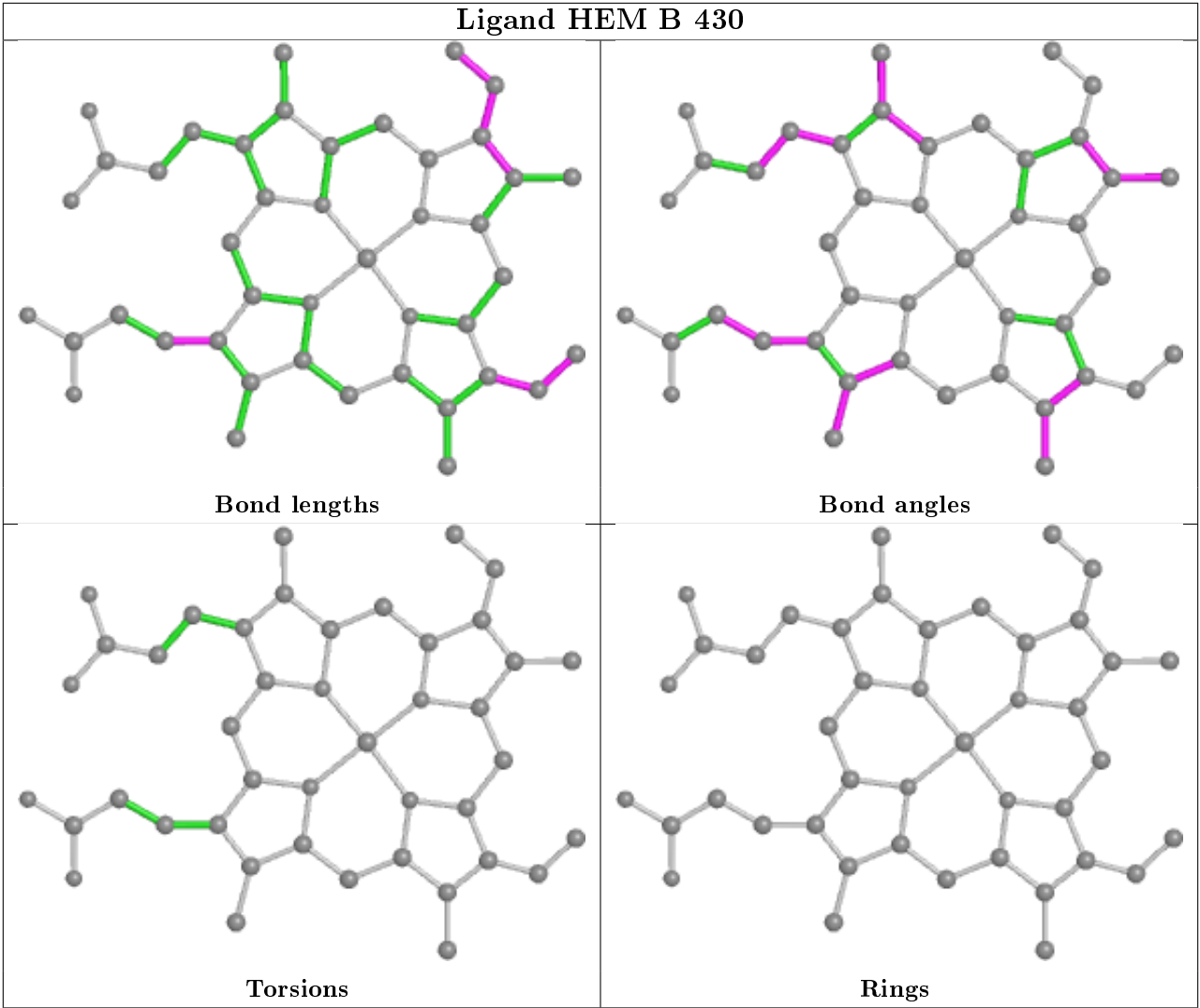
2 monomers are involved in 10 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	632	MES	6	0
2	A	631	SPK	4	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](#)

The following chains have linkage breaks:

Mol	Chain	Number of breaks
1	B	1

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	B	182:SER	C	183:HIS	N	1.19

6 Fit of model and data

6.1 Protein, DNA and RNA chains

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	384/406 (94%)	0.48	37 (9%) 8 10	7, 14, 43, 59	0
1	B	402/406 (99%)	0.51	38 (9%) 8 11	7, 14, 38, 52	0
All	All	786/812 (96%)	0.50	75 (9%) 8 11	7, 14, 41, 59	0

All (75) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	207	SER	15.2
1	B	182	SER	12.6
1	B	87	ILE	10.1
1	A	5	THR	10.1
1	A	204	GLY	9.8
1	B	5	THR	7.7
1	A	88	PRO	7.5
1	A	208	ASP	7.3
1	B	183	HIS	7.3
1	A	207	SER	7.2
1	B	181	SER	7.1
1	B	209	SER	7.1
1	A	172	HIS	7.1
1	A	169	HIS	7.0
1	A	89	ALA	7.0
1	B	180	SER	7.0
1	A	205	LEU	6.8
1	A	87	ILE	6.8
1	A	165	ALA	6.6
1	A	175	THR	6.6
1	B	205	LEU	6.6
1	B	184	GLY	5.6
1	B	187	VAL	5.5
1	A	166	THR	5.4

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Mol	Chain	Res	Type	RSRZ
1	B	98	PRO	5.4
1	B	339	SER	5.3
1	B	208	ASP	5.1
1	B	88	PRO	4.9
1	B	185	ALA	4.9
1	B	90	ARG	4.8
1	A	98	PRO	4.8
1	B	6	ILE	4.7
1	A	164	PRO	4.7
1	B	210	ALA	4.6
1	A	174	TRP	4.5
1	A	86	PHE	4.5
1	B	85	HIS	4.5
1	A	340	PRO	4.4
1	B	140	PRO	4.3
1	A	140	PRO	4.3
1	A	202	LEU	4.1
1	A	195	MET	4.1
1	B	77[A]	ARG	4.0
1	A	206	ARG	3.8
1	B	89	ALA	3.4
1	A	225	ASP	3.3
1	B	91	GLY	3.2
1	B	186	GLU	3.1
1	A	223	GLY	3.1
1	A	168	ARG	3.1
1	A	167	ASP	3.1
1	A	6	ILE	3.1
1	B	340	PRO	3.0
1	A	173	THR	3.0
1	A	163	VAL	2.9
1	B	84	PRO	2.9
1	A	199	PHE	2.8
1	A	194	GLU	2.8
1	A	141	PRO	2.7
1	B	86	PHE	2.7
1	B	141	PRO	2.7
1	B	75	MET	2.7
1	A	339	SER	2.6
1	A	115	ARG	2.5
1	B	83	ALA	2.4
1	A	198	TYR	2.4

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Mol	Chain	Res	Type	RSRZ
1	A	209	SER	2.4
1	A	171	MET	2.4
1	B	99	PRO	2.3
1	B	206	ARG	2.2
1	B	48	GLU	2.2
1	B	179	LEU	2.2
1	B	62	LEU	2.1
1	B	225	ASP	2.1
1	B	212	GLU	2.0

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 carbohydrates in this entry.

6.4 Ligands [i](#)

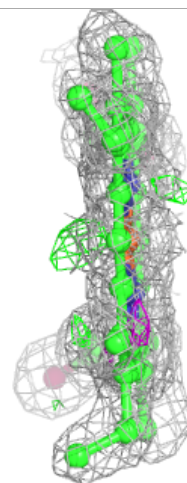
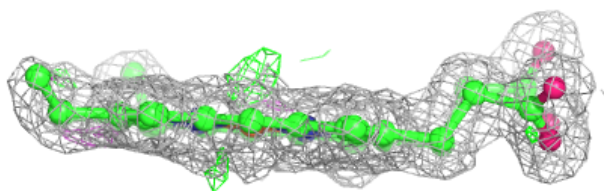
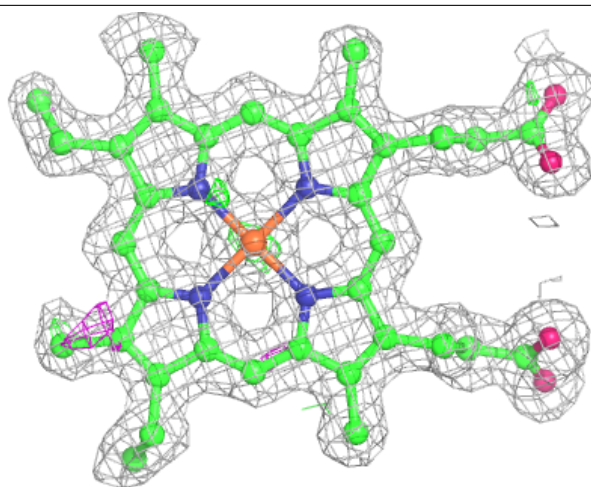
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

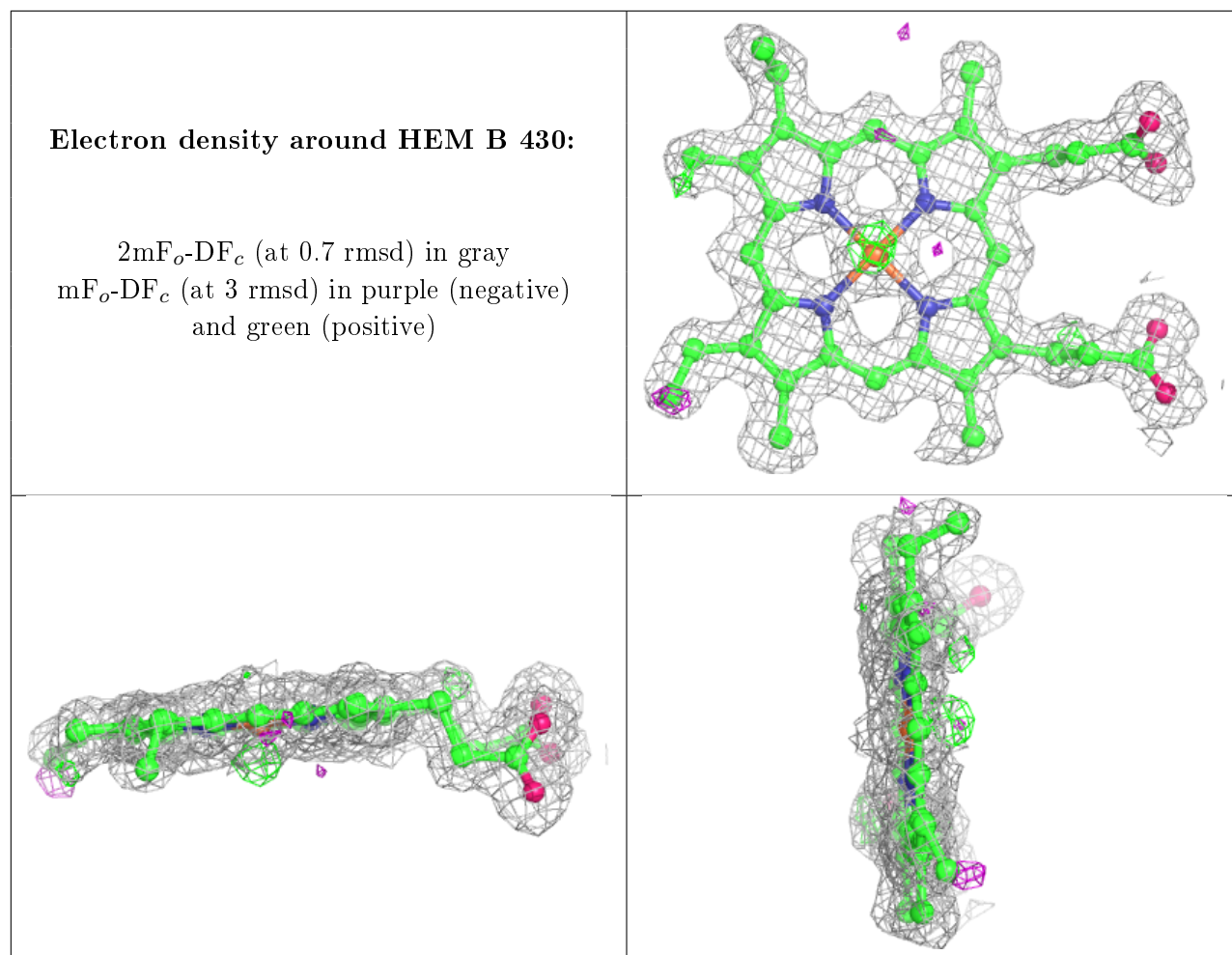
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
4	MES	A	632	12/12	0.33	0.55	47,48,49,50	12
2	SPK	A	631	14/14	0.79	0.20	29,31,33,36	0
3	HEM	A	430	43/43	0.97	0.11	6,9,12,16	0
3	HEM	B	430	43/43	0.98	0.11	6,9,12,16	0

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 430:

$2mF_o-DF_c$ (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.