



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 28, 2022 – 10:26 AM EST

PDB ID : 7SHI
Title : Crystal Structure of Cytochrome P450 AmphL from Streptomyces nodosus and the Structural Basis for Substrate Selectivity in Macrolide Metabolizing P450s
Authors : Amaya, J.A.; Poulos, T.L.
Deposited on : 2021-10-08
Resolution : 2.00 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 : 2.27
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) : 2.27

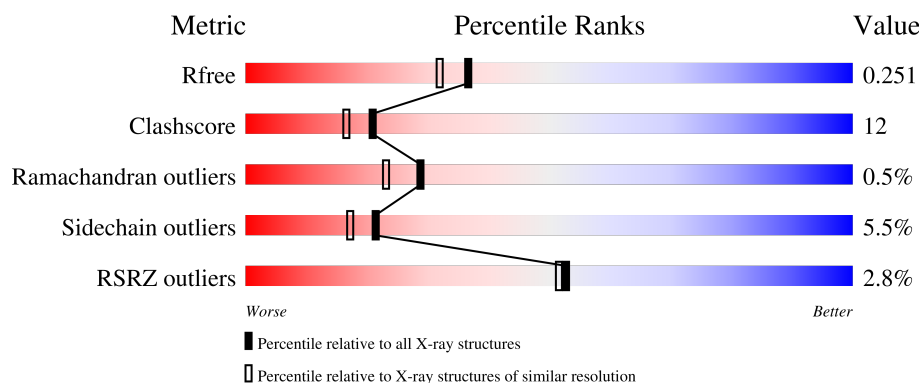
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.00 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	8085 (2.00-2.00)
Clashscore	141614	9178 (2.00-2.00)
Ramachandran outliers	138981	9054 (2.00-2.00)
Sidechain outliers	138945	9053 (2.00-2.00)
RSRZ outliers	127900	7900 (2.00-2.00)

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	391	<div> <div>3%</div> <div>78%</div> <div>19%</div> <div>.</div> </div>
1	B	391	<div> <div>3%</div> <div>78%</div> <div>19%</div> <div>.</div> </div>

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 6798 atoms, of which 150 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 AmphL.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	391	Total	C	N	O	S	0	4	0
			3022	1908	527	573	14			
1	B	391	Total	C	N	O	S	0	3	0
			3017	1903	527	573	14			

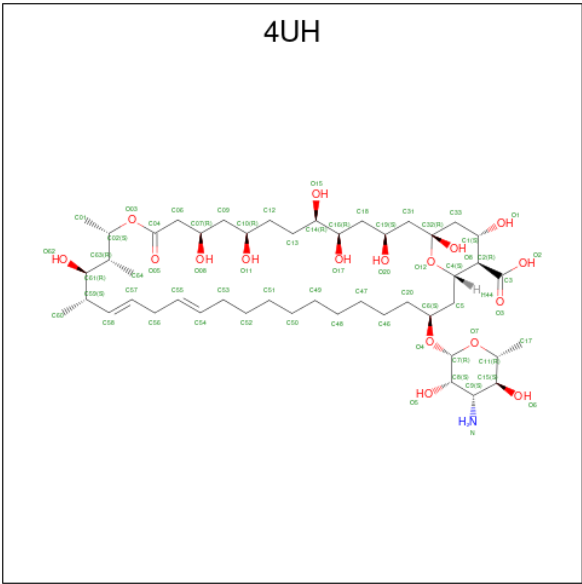
- Molecule 2 is PROTOPORPHYRIN IX CONTAINING FE (three-letter code: HEM) (formula: $C_{34}H_{32}FeN_4O_4$).



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

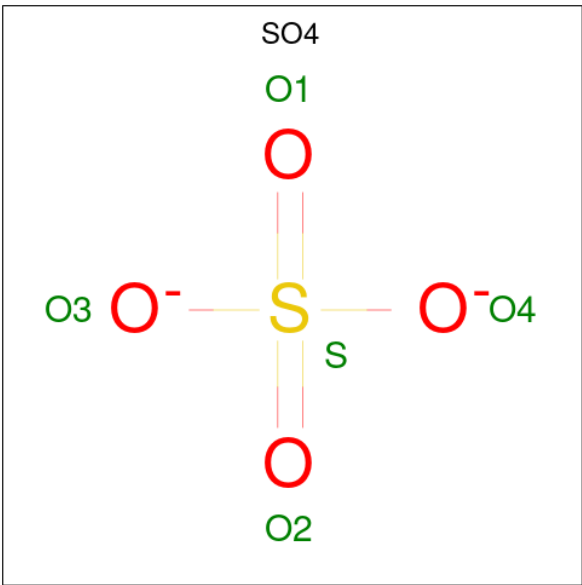
- Molecule 3 is (1R,3S,5R,6R,9R,11R,15S,16R,17R,18S,19E,22E,33S,35S,36R,37S)-33-[(3-amino-3,6-dideoxy-beta-D-mannopyranosyl)oxy]-1,3,5,6,9,11,17,37-octahydroxy-15,16,18-trimethyl-13-oxo-14,39-dioxabicyclo[33.3.1]nonatriaconta-19,22-diene-36-carboxyli

c acid (three-letter code: 4UH) (formula: C₄₇H₈₃NO₁₇) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	H	N	O	0	0
			140	47	75	1	17		
3	B	1	Total	C	H	N	O	0	0
			140	47	75	1	17		

- Molecule 4 is SULFATE ION (three-letter code: SO₄) (formula: O₄S).



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

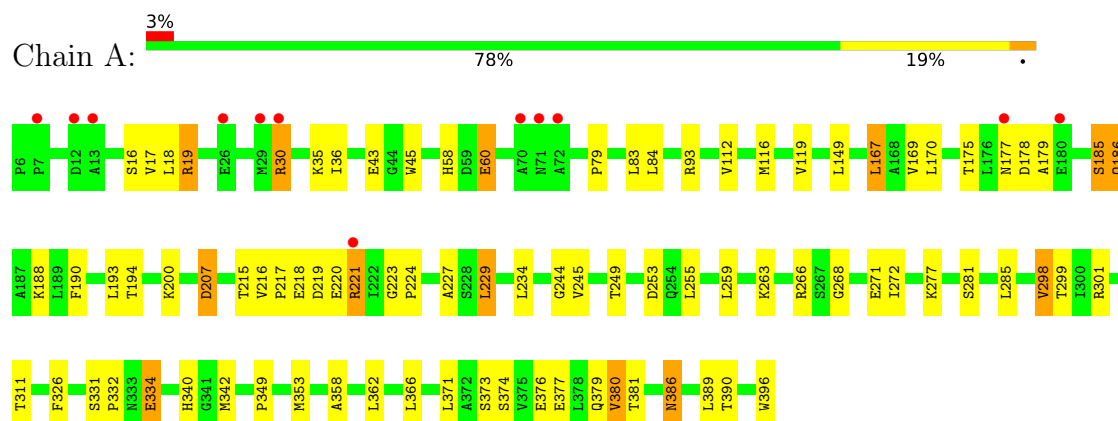
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	180	Total	O	0	0
			180	180		
5	B	198	Total	O	0	0
			198	198		

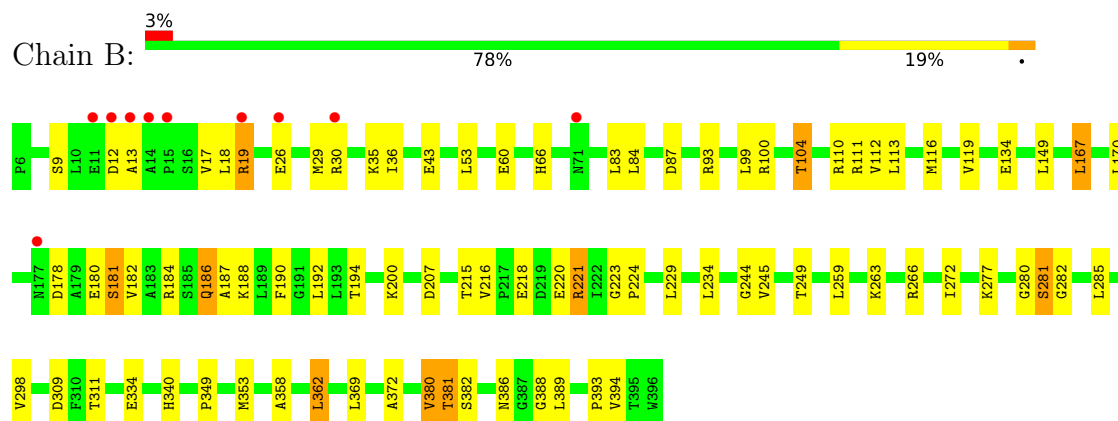
3 Residue-property plots [i](#)

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: AmphL



• Molecule 1: AmphL



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	73.46Å 68.15Å 147.53Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	36.88 – 2.00 36.88 – 2.00	Depositor EDS
% Data completeness (in resolution range)	95.0 (36.88-2.00) 95.0 (36.88-2.00)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.24 (at 2.00Å)	Xtriage
Refinement program	PHENIX 1.11.1 _2575	Depositor
R, R_{free}	0.191 , 0.251 0.191 , 0.251	Depositor DCC
R_{free} test set	1999 reflections (4.14%)	wwPDB-VP
Wilson B-factor (Å ²)	23.0	Xtriage
Anisotropy	0.371	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 52.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	6798	wwPDB-VP
Average B, all atoms (Å ²)	28.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 41.32 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 2.4026e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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: SO4, 4UH, HEM

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z > 5$	RMSZ	$\# Z > 5$
1	A	0.37	0/3079	0.56	0/4186
1	B	0.41	0/3071	0.58	0/4175
All	All	0.39	0/6150	0.57	0/8361

There are no bond length outliers.

There are no bond angle outliers.

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	3022	0	3056	77	0
1	B	3017	0	3045	72	0
2	A	43	0	30	7	0
2	B	43	0	30	4	0
3	A	65	75	0	1	0
3	B	65	75	0	1	0
4	A	10	0	0	0	0
4	B	5	0	0	0	0
5	A	180	0	0	6	0
5	B	198	0	0	5	2
All	All	6648	150	6161	152	2

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 12.

All (152) 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:93:ARG:HH12	1:B:93:ARG:HH12	1.09	0.99
1:B:334:GLU:N	1:B:334:GLU:OE1	2.03	0.90
1:A:175:THR:HG22	1:A:177:ASN:H	1.41	0.85
1:A:194:THR:HG23	1:A:223:GLY:HA3	1.67	0.77
1:A:17:VAL:HG21	1:A:389:LEU:HD21	1.65	0.77
1:A:229:LEU:CD2	2:A:401:HEM:HAC	2.14	0.77
1:A:83:LEU:HD12	1:A:186:GLN:OE1	1.84	0.77
1:A:193:LEU:HD12	1:A:227:ALA:HB2	1.69	0.74
1:B:200:LYS:NZ	5:B:503:HOH:O	2.23	0.72
1:B:170:LEU:HD13	1:B:234:LEU:HD22	1.72	0.71
1:A:58:HIS:ND1	1:B:60:GLU:OE1	2.17	0.71
1:B:186:GLN:HG3	1:B:187:ALA:N	2.04	0.71
1:A:35:LYS:HD2	1:A:43:GLU:OE1	1.91	0.71
1:A:374:SER:OG	1:A:376:GLU:HG2	1.90	0.71
1:B:190:PHE:O	1:B:194:THR:HG23	1.92	0.70
1:B:13:ALA:O	5:B:502:HOH:O	2.11	0.69
1:B:178[B]:ASP:OD2	1:B:181:SER:HB2	1.92	0.69
1:A:200:LYS:NZ	1:A:207:ASP:OD2	2.27	0.68
1:B:178[A]:ASP:OD2	1:B:181:SER:HB2	1.92	0.68
1:A:373:SER:HB2	1:A:377:GLU:OE1	1.93	0.67
1:A:116:MET:HG2	1:A:353:MET:SD	2.34	0.67
1:A:244:GLY:HA2	1:A:272[A]:ILE:HD12	1.77	0.67
1:B:281:SER:OG	1:B:282:GLY:N	2.23	0.67
1:A:178[A]:ASP:OD1	5:A:501:HOH:O	2.13	0.66
1:A:178[B]:ASP:OD1	5:A:501:HOH:O	2.13	0.66
1:A:93:ARG:NH1	1:B:93:ARG:HH12	1.87	0.66
1:B:119:VAL:HG11	1:B:353:MET:HE1	1.78	0.65
1:A:366:LEU:HD22	1:A:396:TRP:CD1	2.32	0.65
1:A:17:VAL:HG12	1:A:18:LEU:HG	1.79	0.65
1:A:19:ARG:HD3	1:A:19:ARG:O	1.96	0.64
1:B:12:ASP:OD2	1:B:382:SER:HB2	1.98	0.64
1:B:194:THR:HG22	1:B:223:GLY:HA3	1.80	0.64
1:A:342:MET:HG3	5:A:536:HOH:O	1.98	0.63
1:B:17:VAL:HG22	1:B:380:VAL:HG12	1.81	0.63
1:B:358:ALA:O	1:B:362:LEU:HB2	1.99	0.62
1:B:192:LEU:HD12	1:B:192:LEU:O	1.99	0.62
1:A:17:VAL:HG21	1:A:389:LEU:CD2	2.30	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:299:THR:HG22	1:A:301:ARG:HD3	1.82	0.61
1:B:263:LYS:HD3	1:B:266:ARG:HE	1.65	0.61
1:A:218:GLU:N	1:A:218:GLU:OE2	2.33	0.61
1:A:60:GLU:O	1:A:60:GLU:HG2	2.00	0.61
1:B:119:VAL:HG11	1:B:353:MET:CE	2.31	0.60
1:A:194:THR:CG2	1:A:223:GLY:HA3	2.30	0.60
1:B:244:GLY:HA2	1:B:272:ILE:HD12	1.82	0.60
1:B:263:LYS:HE3	1:B:266:ARG:HH21	1.67	0.60
1:B:381:THR:HG23	1:B:388:GLY:O	2.01	0.60
1:A:93:ARG:HH12	1:B:93:ARG:NH1	1.91	0.59
1:A:175:THR:CG2	1:A:177:ASN:H	2.11	0.59
1:B:277:LYS:HG3	1:B:311:THR:HG23	1.84	0.59
1:B:216:VAL:HG13	1:B:221:ARG:HD3	1.83	0.59
1:A:221:ARG:CZ	1:A:221:ARG:HB2	2.32	0.58
1:B:35:LYS:HD2	1:B:43:GLU:OE1	2.04	0.58
1:A:119:VAL:HB	1:A:353:MET:HE1	1.85	0.57
1:A:218:GLU:HB3	1:A:220:GLU:OE1	2.04	0.57
1:A:285:LEU:HD21	3:A:402:4UH:C50	2.35	0.57
1:A:30[B]:ARG:NH1	1:A:30[B]:ARG:HA	2.19	0.57
1:B:26:GLU:OE1	1:B:30[B]:ARG:HG2	2.05	0.57
1:A:379:GLN:O	1:A:390[A]:THR:HG22	2.05	0.56
1:B:12:ASP:CG	1:B:382:SER:HB2	2.25	0.56
1:B:17:VAL:CG2	1:B:380:VAL:HG12	2.35	0.56
1:B:19:ARG:O	1:B:19:ARG:HD3	2.05	0.56
1:B:334:GLU:H	1:B:334:GLU:CD	2.07	0.55
1:A:229:LEU:HD22	2:A:401:HEM:HAC	1.89	0.55
1:A:358:ALA:O	1:A:362:LEU:HB2	2.07	0.54
1:B:372:ALA:HB3	1:B:393:PRO:HB2	1.88	0.54
1:A:16:SER:OG	1:A:19:ARG:HD2	2.08	0.54
1:A:167:LEU:O	5:A:502:HOH:O	2.19	0.54
1:A:334:GLU:H	1:A:334:GLU:CD	2.10	0.54
1:B:263:LYS:CD	1:B:266:ARG:HH21	2.21	0.53
1:B:229:LEU:HD22	2:B:401:HEM:HAC	1.91	0.53
1:A:45:TRP:CZ3	1:A:298:VAL:HG21	2.44	0.52
1:B:17:VAL:HG21	1:B:389:LEU:HD21	1.91	0.52
1:A:79:PRO:HG2	1:A:179:ALA:O	2.10	0.52
1:A:112:VAL:HG21	1:A:349:PRO:HB2	1.90	0.52
1:A:84:LEU:HD21	1:A:285:LEU:HD13	1.92	0.51
1:A:149:LEU:HD22	1:A:167:LEU:HD13	1.92	0.51
1:A:331:SER:HA	1:A:332:PRO:C	2.30	0.51
1:B:53:LEU:C	1:B:53:LEU:HD23	2.31	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:17:VAL:HG12	1:B:18:LEU:HG	1.93	0.51
1:A:119:VAL:CB	1:A:353:MET:HE1	2.41	0.50
1:B:190:PHE:CE2	1:B:224:PRO:HB3	2.46	0.50
1:A:93:ARG:NH2	5:A:515:HOH:O	2.44	0.50
1:A:194:THR:HG22	1:A:219:ASP:O	2.11	0.49
1:B:53:LEU:HD23	1:B:53:LEU:O	2.12	0.49
1:A:229:LEU:HD23	2:A:401:HEM:HAC	1.92	0.49
1:B:216:VAL:HG11	1:B:221:ARG:HG3	1.95	0.49
1:B:282:GLY:HA3	1:B:309:ASP:OD1	2.13	0.49
1:B:180:GLU:O	1:B:184:ARG:HG3	2.12	0.49
1:B:285:LEU:HD21	3:B:402:4UH:C50	2.43	0.49
1:A:190:PHE:CE2	1:A:224:PRO:HB3	2.48	0.49
1:A:263:LYS:HD3	1:A:266:ARG:HH21	1.78	0.49
1:B:93:ARG:NH2	5:B:513:HOH:O	2.44	0.49
1:B:178[A]:ASP:O	1:B:182:VAL:HG23	2.12	0.49
1:B:178[B]:ASP:O	1:B:182:VAL:HG23	2.12	0.49
1:B:116:MET:HG2	1:B:353:MET:SD	2.53	0.48
1:B:119:VAL:HB	1:B:353:MET:HE3	1.95	0.48
1:B:263:LYS:CE	1:B:266:ARG:HH21	2.25	0.48
1:B:83:LEU:HD11	1:B:186:GLN:OE1	2.13	0.48
1:B:263:LYS:HE3	1:B:266:ARG:NH2	2.29	0.47
1:A:19:ARG:NH2	5:A:514:HOH:O	2.43	0.47
1:A:216:VAL:HG11	1:A:221:ARG:HG3	1.95	0.47
1:B:60:GLU:HG2	1:B:60:GLU:O	2.15	0.47
1:A:217:PRO:HG2	1:A:218:GLU:OE2	2.15	0.47
1:A:17:VAL:CG1	1:A:18:LEU:HG	2.44	0.47
1:A:119:VAL:HG11	1:A:353:MET:CE	2.45	0.47
1:B:26:GLU:OE1	1:B:30[B]:ARG:NE	2.47	0.47
1:A:30[B]:ARG:HA	1:A:30[B]:ARG:HH11	1.78	0.46
1:A:220:GLU:H	1:A:220:GLU:CD	2.19	0.46
1:B:83:LEU:CD1	1:B:186:GLN:OE1	2.63	0.46
1:A:36:ILE:C	1:A:36:ILE:HD12	2.36	0.46
1:B:36:ILE:HD12	1:B:36:ILE:C	2.36	0.46
1:B:111:ARG:NH2	5:B:516:HOH:O	2.49	0.46
1:B:26:GLU:OE2	1:B:29:MET:HE3	2.14	0.45
1:A:379:GLN:HB2	1:A:390[A]:THR:CG2	2.47	0.45
1:B:149:LEU:HD22	1:B:167:LEU:HD13	1.98	0.45
1:B:13:ALA:HB1	1:B:380:VAL:HG21	1.99	0.45
2:B:401:HEM:HMB2	2:B:401:HEM:HBB2	1.99	0.45
1:A:271:GLU:HG3	1:A:326:PHE:CE1	2.51	0.44
1:B:112:VAL:HG21	1:B:349:PRO:HB2	1.98	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:223:GLY:N	1:A:224:PRO:HD2	2.33	0.44
2:B:401:HEM:HBB2	2:B:401:HEM:CMB	2.48	0.44
1:A:271:GLU:HG3	1:A:326:PHE:CD1	2.53	0.44
1:B:66:HIS:O	1:B:87:ASP:HA	2.17	0.44
2:A:401:HEM:CMC	2:A:401:HEM:HBC2	2.48	0.44
1:A:193:LEU:CD1	1:A:227:ALA:HB2	2.45	0.43
1:B:134:GLU:HA	1:B:134:GLU:OE1	2.17	0.43
1:A:83:LEU:HD23	1:A:83:LEU:HA	1.79	0.43
1:A:169:VAL:HG13	1:A:185:SER:HB3	2.00	0.43
1:A:268:GLY:O	1:A:272[A]:ILE:HG12	2.18	0.43
1:A:175:THR:HG22	1:A:177:ASN:N	2.20	0.43
1:B:280:GLY:O	1:B:281:SER:C	2.57	0.43
1:A:17:VAL:HG23	1:A:380:VAL:HG12	1.99	0.42
1:B:113:LEU:HD23	5:B:615:HOH:O	2.18	0.42
1:A:245:VAL:O	1:A:249:THR:HG23	2.18	0.42
2:B:401:HEM:HBC2	2:B:401:HEM:HMC2	2.01	0.42
1:A:170:LEU:HD13	1:A:234:LEU:HD22	2.01	0.41
1:B:221:ARG:CZ	1:B:221:ARG:HB2	2.49	0.41
1:A:249:THR:OG1	1:A:371:LEU:HD21	2.20	0.41
1:A:249:THR:OG1	1:A:371:LEU:CD2	2.68	0.41
2:A:401:HEM:HBC2	2:A:401:HEM:HMC2	2.02	0.41
2:A:401:HEM:HBB2	2:A:401:HEM:HMB2	2.01	0.41
1:B:188:LYS:HE2	1:B:188:LYS:HB3	1.72	0.41
1:B:218:GLU:HB3	1:B:220:GLU:OE1	2.21	0.41
1:A:277:LYS:HG3	1:A:311:THR:HG23	2.02	0.41
1:B:245:VAL:O	1:B:249:THR:HG23	2.21	0.41
1:A:188:LYS:HE2	1:A:188:LYS:HB3	1.65	0.41
1:A:255:LEU:HD22	1:A:255:LEU:O	2.21	0.40
1:B:100:ARG:O	1:B:104:THR:HB	2.20	0.40
1:B:263:LYS:HD3	1:B:263:LYS:HA	1.85	0.40
2:A:401:HEM:HBB2	2:A:401:HEM:CMB	2.51	0.40
1:B:369:LEU:HD11	1:B:394:VAL:HG21	2.02	0.40
1:A:253:ASP:OD1	1:A:253:ASP:N	2.55	0.40

All (2) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:B:566:HOH:O	5:B:669:HOH:O[3_754]	2.13	0.07
5:B:616:HOH:O	5:B:644:HOH:O[3_744]	2.15	0.05

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	393/391 (100%)	381 (97%)	10 (2%)	2 (0%)	29	23
1	B	392/391 (100%)	382 (97%)	8 (2%)	2 (0%)	29	23
All	All	785/782 (100%)	763 (97%)	18 (2%)	4 (0%)	29	23

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	281	SER
1	A	386	ASN
1	B	281	SER
1	B	386	ASN

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	322/318 (101%)	304 (94%)	18 (6%)	21	17
1	B	321/318 (101%)	303 (94%)	18 (6%)	21	17
All	All	643/636 (101%)	607 (94%)	36 (6%)	21	17

All (36) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	19	ARG
1	A	30[A]	ARG

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Mol	Chain	Res	Type
1	A	30[B]	ARG
1	A	60	GLU
1	A	167	LEU
1	A	185	SER
1	A	186	GLN
1	A	207	ASP
1	A	215	THR
1	A	221	ARG
1	A	229	LEU
1	A	259	LEU
1	A	298	VAL
1	A	334	GLU
1	A	340	HIS
1	A	380	VAL
1	A	381	THR
1	A	386	ASN
1	B	9	SER
1	B	19	ARG
1	B	84	LEU
1	B	99	LEU
1	B	104	THR
1	B	110	ARG
1	B	167	LEU
1	B	181	SER
1	B	186	GLN
1	B	207	ASP
1	B	215	THR
1	B	221	ARG
1	B	259	LEU
1	B	298	VAL
1	B	340	HIS
1	B	362	LEU
1	B	380	VAL
1	B	381	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no monosaccharides in this entry.

5.6 Ligand geometry [i](#)

7 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$
2	HEM	B	401	1	27,50,50	1.91	5 (18%)	17,82,82	1.65	3 (17%)
3	4UH	A	402	-	62,67,67	1.34	7 (11%)	76,92,92	1.81	17 (22%)
4	SO4	A	403	-	4,4,4	0.19	0	6,6,6	0.16	0
4	SO4	B	403	-	4,4,4	0.17	0	6,6,6	0.50	0
4	SO4	A	404	-	4,4,4	0.10	0	6,6,6	0.26	0
2	HEM	A	401	1	27,50,50	1.65	4 (14%)	17,82,82	1.71	4 (23%)
3	4UH	B	402	-	62,67,67	1.35	8 (12%)	76,92,92	1.92	17 (22%)

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	401	1	-	0/6/54/54	-
2	HEM	A	401	1	-	0/6/54/54	-
3	4UH	B	402	-	-	23/67/109/109	0/2/3/3
3	4UH	A	402	-	-	25/67/109/109	0/2/3/3

All (24) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	401	HEM	C3B-C2B	-4.66	1.33	1.40
2	B	401	HEM	C3C-C2C	-4.40	1.34	1.40
2	B	401	HEM	C3B-CAB	3.64	1.55	1.47
2	A	401	HEM	C3C-C2C	-3.47	1.35	1.40
3	A	402	4UH	C50-C49	-3.38	1.32	1.51
2	A	401	HEM	C3B-C2B	-3.35	1.35	1.40
2	A	401	HEM	C3B-CAB	3.35	1.54	1.47
3	B	402	4UH	C48-C47	-3.24	1.33	1.51
3	A	402	4UH	C52-C51	-3.22	1.33	1.51
2	B	401	HEM	C3C-CAC	3.20	1.54	1.47
2	A	401	HEM	C3C-CAC	3.18	1.54	1.47
3	A	402	4UH	C48-C47	-3.18	1.33	1.51
3	B	402	4UH	C50-C49	-3.15	1.33	1.51
3	B	402	4UH	C52-C51	-3.01	1.34	1.51
3	A	402	4UH	C53-C54	-2.80	1.34	1.50
3	B	402	4UH	C53-C54	-2.73	1.34	1.50
3	B	402	4UH	C9-N	2.43	1.51	1.47
3	A	402	4UH	C9-N	2.43	1.51	1.47
2	B	401	HEM	CAA-C2A	2.34	1.55	1.52
3	B	402	4UH	C55-C54	2.34	1.45	1.31
3	B	402	4UH	C56-C55	-2.33	1.34	1.51
3	A	402	4UH	C56-C55	-2.29	1.34	1.51
3	A	402	4UH	C55-C54	2.21	1.44	1.31
3	B	402	4UH	C64-C63	2.09	1.57	1.53

All (41) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	402	4UH	C7-O4-C6	7.70	130.61	114.88
3	B	402	4UH	C7-O4-C6	7.38	129.96	114.88
3	B	402	4UH	O7-C11-C15	-5.73	99.25	109.52
3	B	402	4UH	O4-C7-C8	4.37	119.43	108.10
3	A	402	4UH	C02-O03-C04	3.93	125.62	118.55
3	A	402	4UH	O4-C7-C8	3.83	118.03	108.10
3	B	402	4UH	C64-C63-C61	3.75	118.98	111.39
3	A	402	4UH	O7-C11-C15	-3.74	102.81	109.52
3	A	402	4UH	C64-C63-C61	3.73	118.94	111.39
3	B	402	4UH	O03-C02-C63	3.53	115.43	106.30
3	A	402	4UH	O03-C02-C63	3.50	115.36	106.30
2	B	401	HEM	CMD-C2D-C1D	-3.35	123.31	128.46
3	B	402	4UH	C02-O03-C04	3.34	124.55	118.55
3	A	402	4UH	O4-C7-O7	-3.07	102.09	110.67
3	B	402	4UH	O62-C61-C63	-3.00	102.82	109.49

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	401	HEM	CBA-CAA-C2A	-2.99	106.98	112.49
3	B	402	4UH	C11-C15-C9	-2.87	106.16	110.51
3	B	402	4UH	O4-C6-C5	2.86	113.41	108.18
2	A	401	HEM	CMA-C3A-C4A	-2.77	124.20	128.46
3	B	402	4UH	C47-C46-C20	-2.75	103.90	113.62
3	A	402	4UH	C7-O7-C11	2.72	118.35	113.67
3	B	402	4UH	O4-C7-O7	-2.71	103.10	110.67
3	A	402	4UH	C59-C61-C63	2.71	120.49	114.96
3	B	402	4UH	C17-C11-C15	2.69	118.03	113.07
3	B	402	4UH	C57-C56-C55	2.62	124.92	112.02
3	A	402	4UH	O5-C8-C9	-2.61	105.54	110.22
2	A	401	HEM	CMB-C2B-C3B	2.51	129.38	124.68
3	A	402	4UH	O05-C04-C06	-2.50	119.20	124.73
3	A	402	4UH	C57-C56-C55	2.38	123.74	112.02
3	A	402	4UH	C50-C49-C48	2.34	126.29	114.42
2	A	401	HEM	CBD-CAD-C3D	-2.32	108.20	112.48
3	A	402	4UH	O62-C61-C63	-2.30	104.37	109.49
3	A	402	4UH	C47-C46-C20	-2.27	105.60	113.62
3	B	402	4UH	C50-C49-C48	2.19	125.52	114.42
3	A	402	4UH	C11-C15-C9	-2.16	107.24	110.51
2	A	401	HEM	CMC-C2C-C3C	2.12	128.64	124.68
3	A	402	4UH	C49-C48-C47	2.07	124.96	114.42
3	B	402	4UH	C52-C53-C54	2.07	124.32	112.43
3	B	402	4UH	O5-C8-C7	2.04	114.99	110.05
2	B	401	HEM	C4A-C3A-C2A	2.03	108.41	107.00
3	B	402	4UH	O8-C32-C31	-2.01	108.23	111.62

There are no chirality outliers.

All (48) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	402	4UH	C46-C20-C6-O4
3	A	402	4UH	C63-C02-O03-C04
3	A	402	4UH	C59-C61-C63-C64
3	A	402	4UH	O62-C61-C63-C64
3	A	402	4UH	C8-C7-O4-C6
3	B	402	4UH	C46-C20-C6-O4
3	B	402	4UH	C63-C02-O03-C04
3	B	402	4UH	C59-C61-C63-C64
3	B	402	4UH	O62-C61-C63-C64
3	B	402	4UH	C8-C7-O4-C6
3	A	402	4UH	O62-C61-C63-C02

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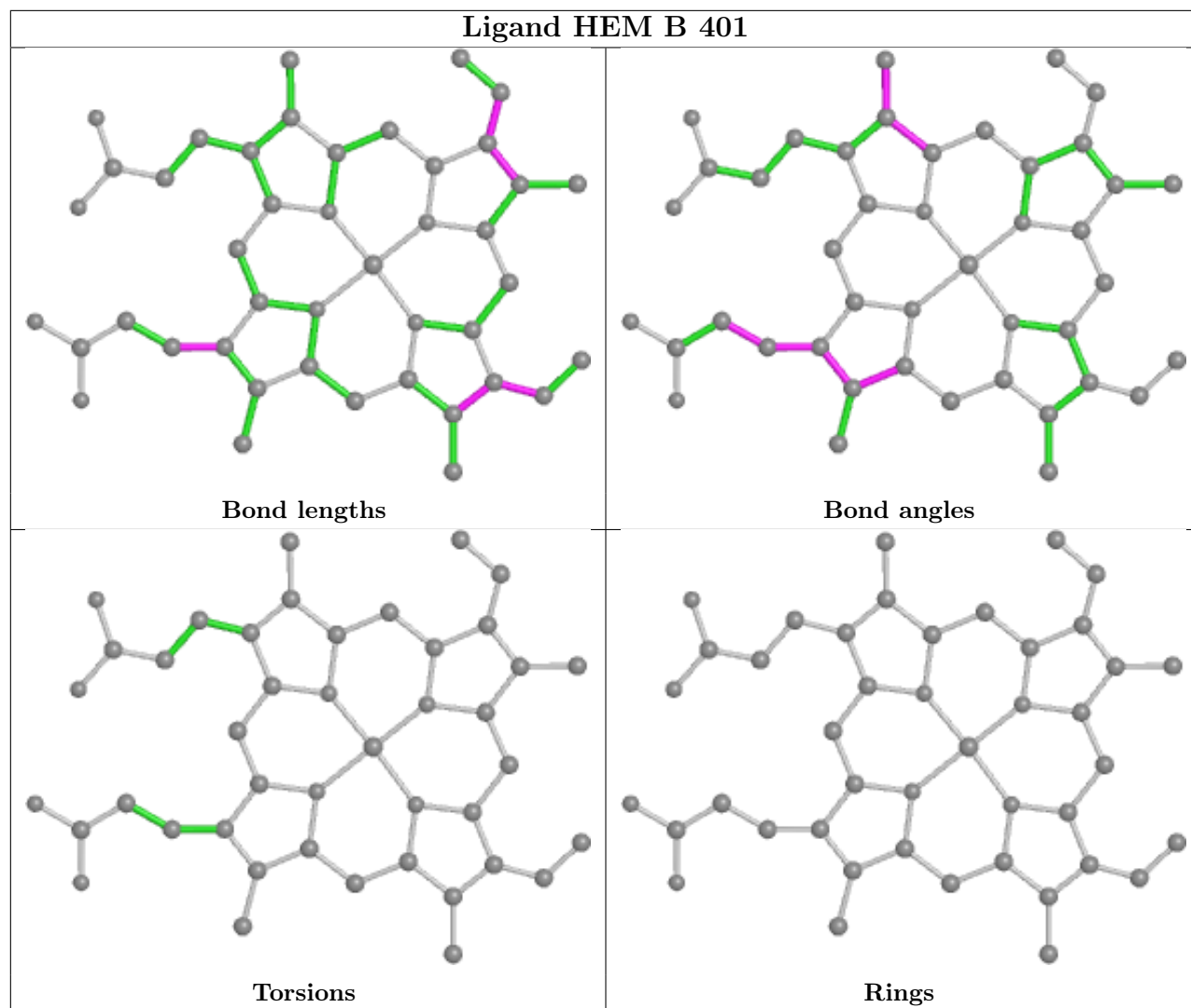
Mol	Chain	Res	Type	Atoms
3	B	402	4UH	O62-C61-C63-C02
3	A	402	4UH	C07-C09-C10-O11
3	B	402	4UH	C59-C61-C63-C02
3	A	402	4UH	C20-C46-C47-C48
3	A	402	4UH	O05-C04-O03-C02
3	B	402	4UH	C47-C48-C49-C50
3	A	402	4UH	C01-C02-C63-C64
3	B	402	4UH	C01-C02-C63-C64
3	B	402	4UH	C49-C50-C51-C52
3	A	402	4UH	C47-C48-C49-C50
3	B	402	4UH	C48-C49-C50-C51
3	A	402	4UH	C59-C61-C63-C02
3	A	402	4UH	C50-C51-C52-C53
3	B	402	4UH	O05-C04-O03-C02
3	B	402	4UH	C06-C04-O03-C02
3	B	402	4UH	C46-C47-C48-C49
3	A	402	4UH	C49-C50-C51-C52
3	A	402	4UH	C06-C04-O03-C02
3	B	402	4UH	C20-C46-C47-C48
3	B	402	4UH	C50-C51-C52-C53
3	A	402	4UH	C54-C55-C56-C57
3	B	402	4UH	C54-C55-C56-C57
3	A	402	4UH	C46-C20-C6-C5
3	A	402	4UH	C46-C47-C48-C49
3	B	402	4UH	C57-C58-C59-C60
3	A	402	4UH	C51-C52-C53-C54
3	A	402	4UH	C57-C58-C59-C60
3	A	402	4UH	C55-C56-C57-C58
3	B	402	4UH	C55-C56-C57-C58
3	A	402	4UH	C01-C02-O03-C04
3	B	402	4UH	C01-C02-O03-C04
3	A	402	4UH	C48-C49-C50-C51
3	B	402	4UH	C52-C53-C54-C55
3	A	402	4UH	C52-C53-C54-C55
3	B	402	4UH	C60-C59-C61-C63
3	A	402	4UH	O03-C02-C63-C64
3	B	402	4UH	O03-C02-C63-C64

There are no ring outliers.

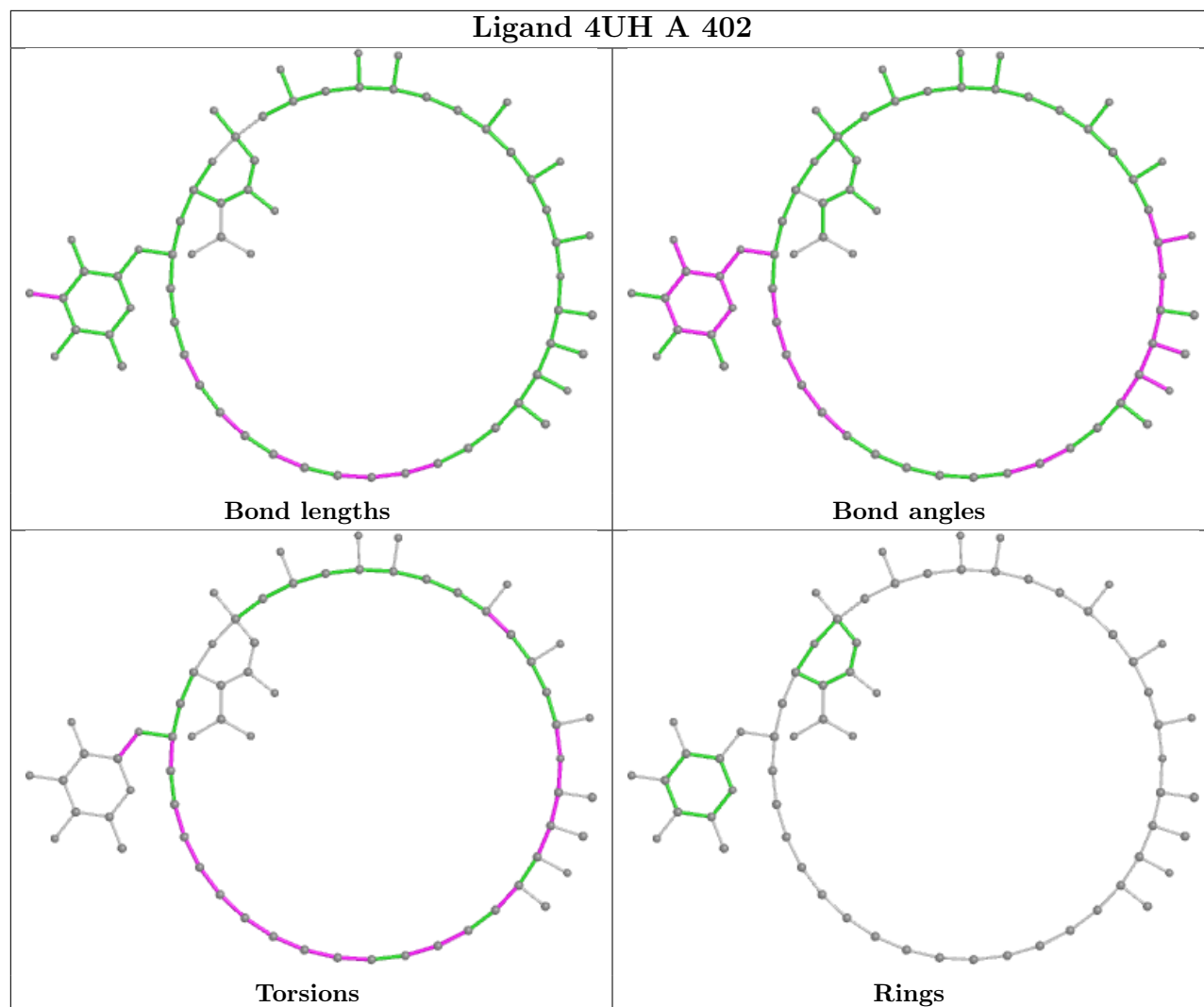
4 monomers are involved in 13 short contacts:

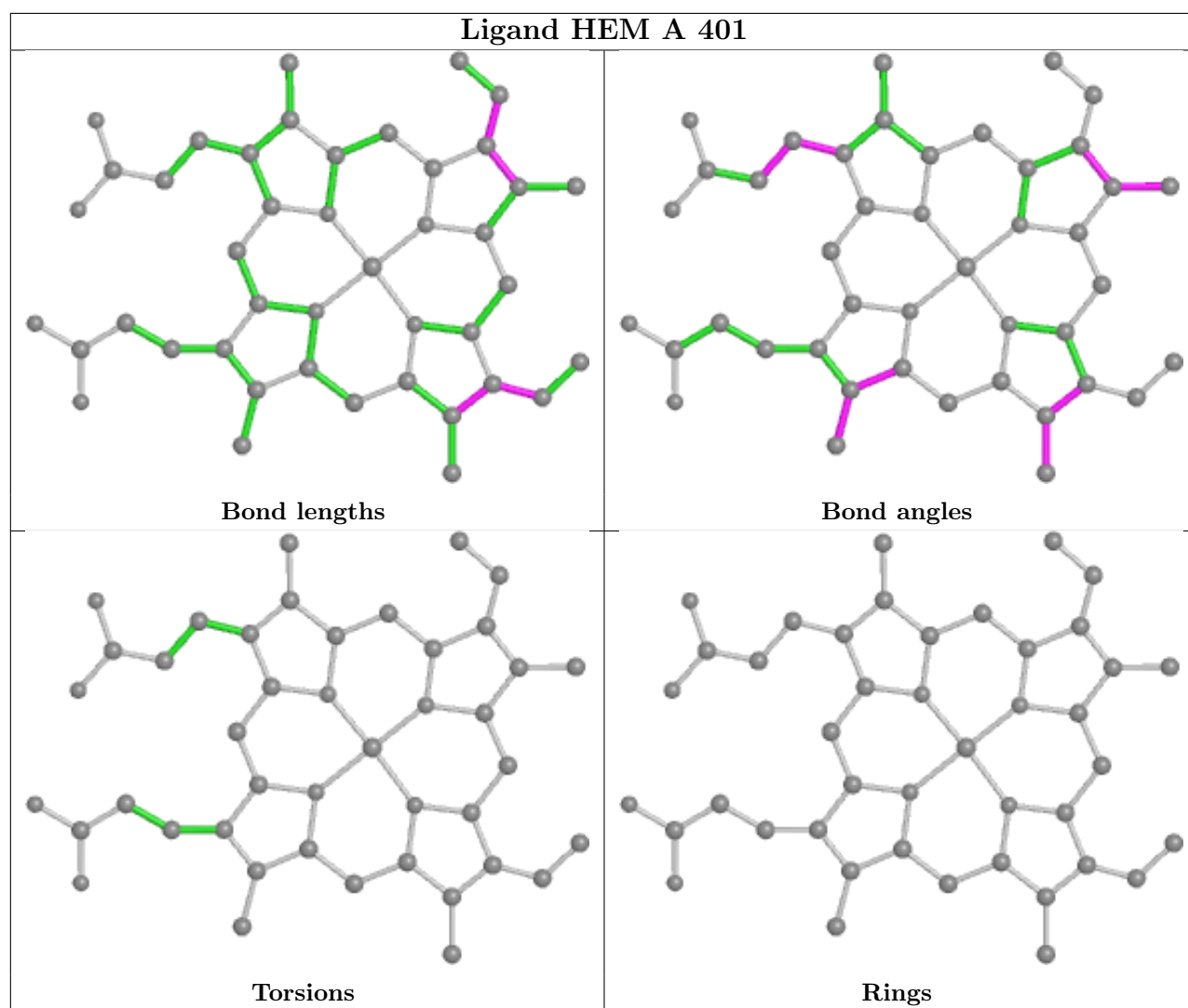
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	401	HEM	4	0
3	A	402	4UH	1	0
2	A	401	HEM	7	0
3	B	402	4UH	1	0

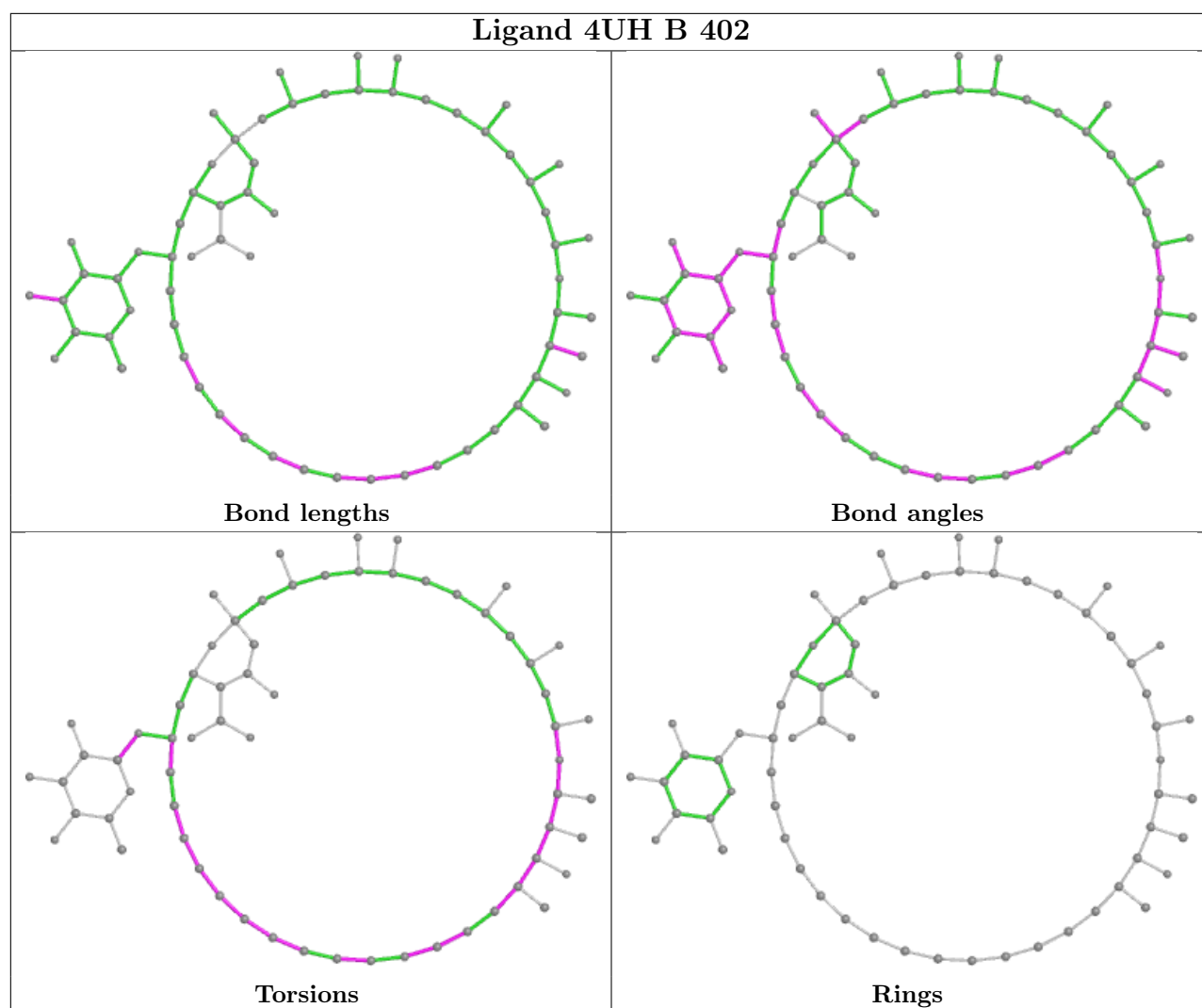
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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



Ligand 4UH A 402







5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

In the following table, the column labelled ‘#RSRZ > 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 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	391/391 (100%)	0.02	12 (3%)	49	48	13, 27, 47, 68	0
1	B	391/391 (100%)	-0.03	10 (2%)	56	54	12, 24, 43, 84	0
All	All	782/782 (100%)	-0.00	22 (2%)	53	51	12, 25, 45, 84	0

All (22) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	14	ALA	10.9
1	B	13	ALA	6.2
1	B	12	ASP	4.9
1	B	15	PRO	4.3
1	A	30[A]	ARG	4.1
1	B	177	ASN	3.5
1	A	70	ALA	3.2
1	A	12	ASP	3.0
1	A	13	ALA	2.8
1	A	180	GLU	2.6
1	B	11	GLU	2.6
1	A	7	PRO	2.4
1	B	26	GLU	2.4
1	A	29	MET	2.3
1	B	71	ASN	2.3
1	A	72	ALA	2.3
1	A	26	GLU	2.2
1	B	19	ARG	2.2
1	A	221	ARG	2.2
1	A	71	ASN	2.1
1	B	30[A]	ARG	2.0
1	A	177	ASN	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 monosaccharides in this entry.

6.4 Ligands [i](#)

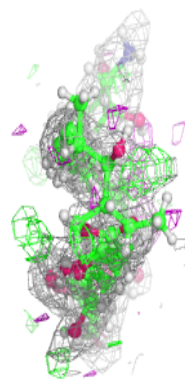
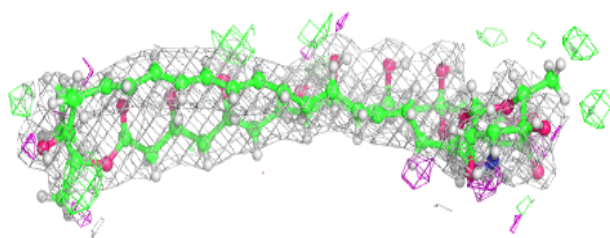
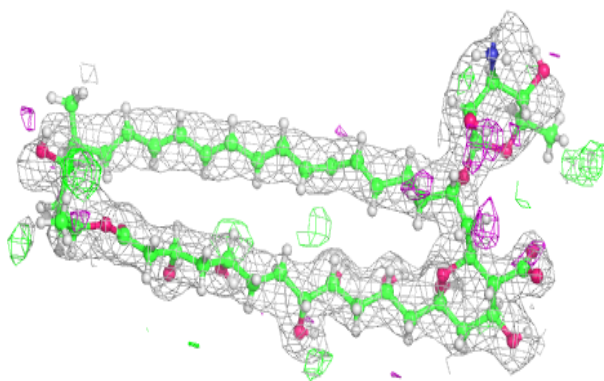
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
3	4UH	B	402	65/65	0.90	0.17	15,29,63,80	0
3	4UH	A	402	65/65	0.91	0.16	16,31,66,69	0
4	SO4	A	404	5/5	0.96	0.10	31,34,37,38	5
2	HEM	B	401	43/43	0.98	0.14	8,13,19,23	0
4	SO4	A	403	5/5	0.99	0.08	25,25,28,37	0
2	HEM	A	401	43/43	0.99	0.13	9,12,21,25	0
4	SO4	B	403	5/5	0.99	0.12	25,26,28,30	0

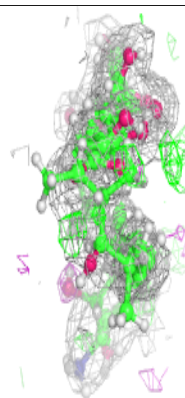
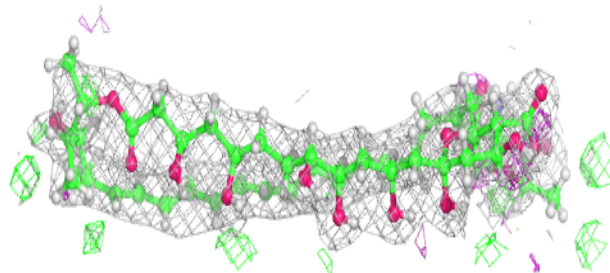
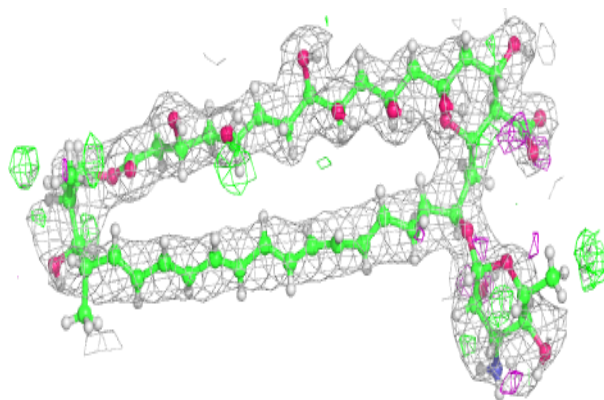
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around 4UH B 402:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

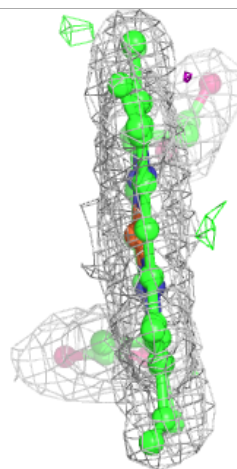
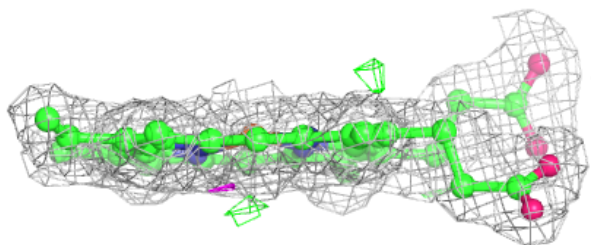
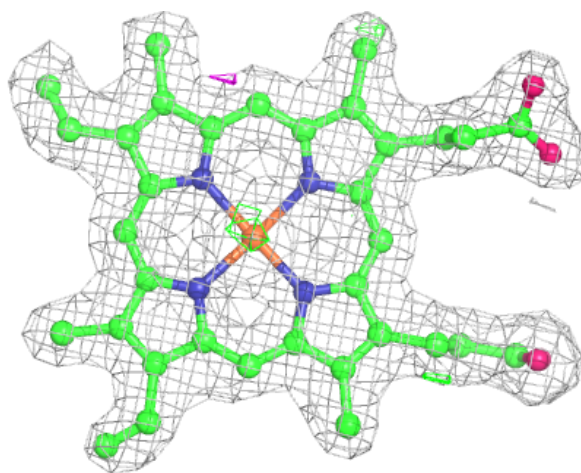
**Electron density around 4UH A 402:**

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



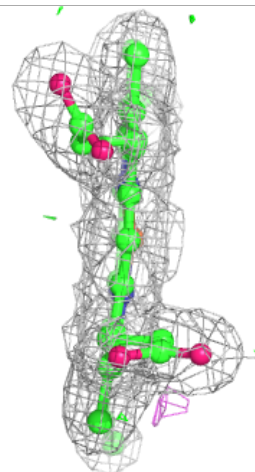
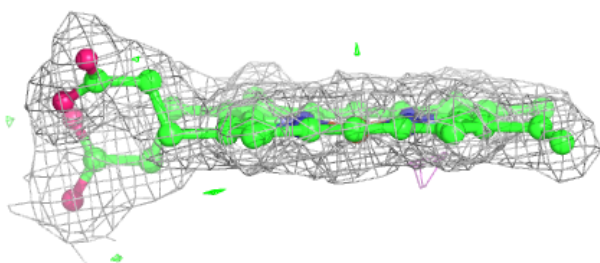
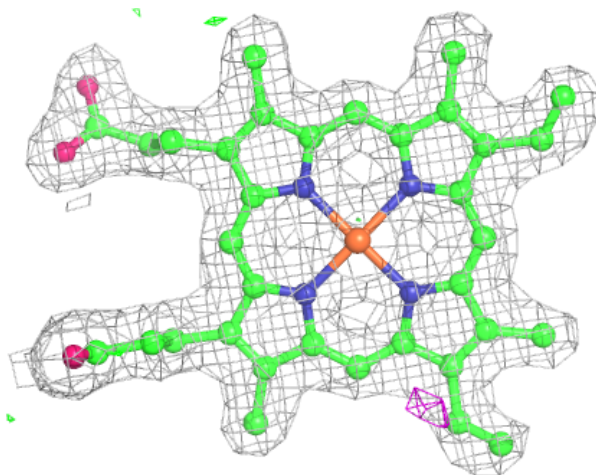
Electron density around HEM B 401:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around HEM A 401:

$2mF_o-DF_c$ (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.