



Full wwPDB X-ray Structure Validation Report ⓘ

Aug 4, 2026 – 02:00 PM EDT

PDB ID : 1S1F / pdb_00001s1f
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 : 2004-01-06
Resolution : 1.50 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.015 (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

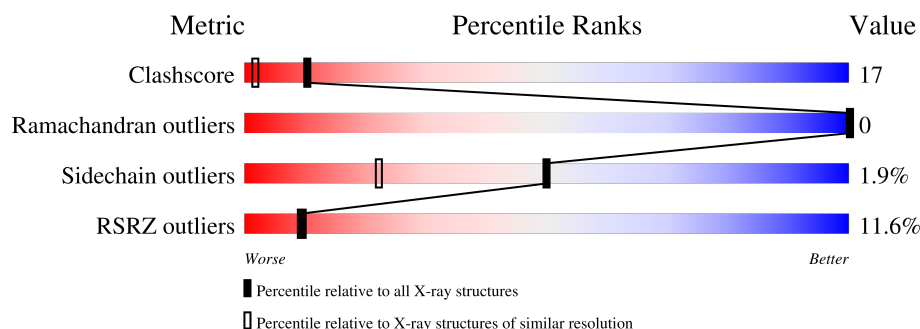
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.50 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))
Clashscore	190562	4235 (1.50-1.50)
Ramachandran outliers	187476	4153 (1.50-1.50)
Sidechain outliers	187428	4150 (1.50-1.50)
RSRZ outliers	180081	4039 (1.50-1.50)

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	406	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
5	MLA	A	433	-	X	-	-
6	GOL	A	434	-	X	-	-

2 Entry composition [i](#)

There are 7 unique types of molecules in this entry. The entry contains 3495 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	397	Total	C	N	O	S	0	4	0
			3093	1946	569	567	11			

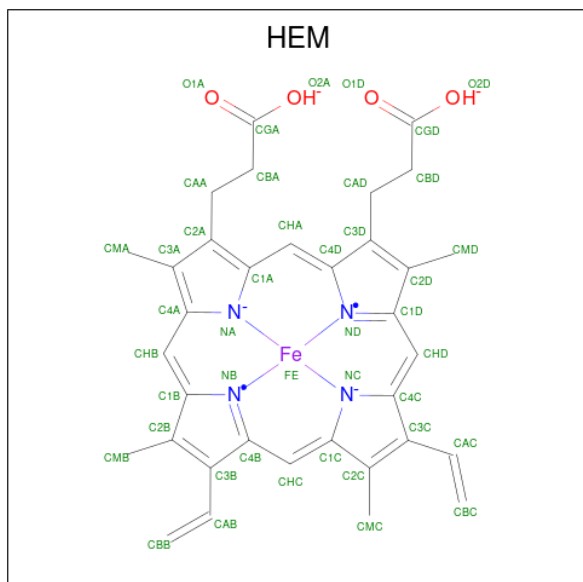
There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	405	HIS	-	expression tag	GB 21219717
A	406	ALA	-	expression tag	GB 21219717

- Molecule 2 is MERCURY (II) ION (CCD ID: HG) (formula: Hg).

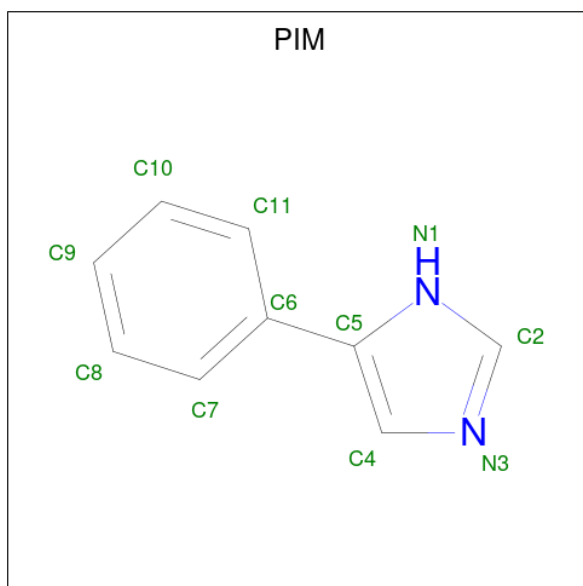
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	2	Total	Hg	0	0
			2	2		

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



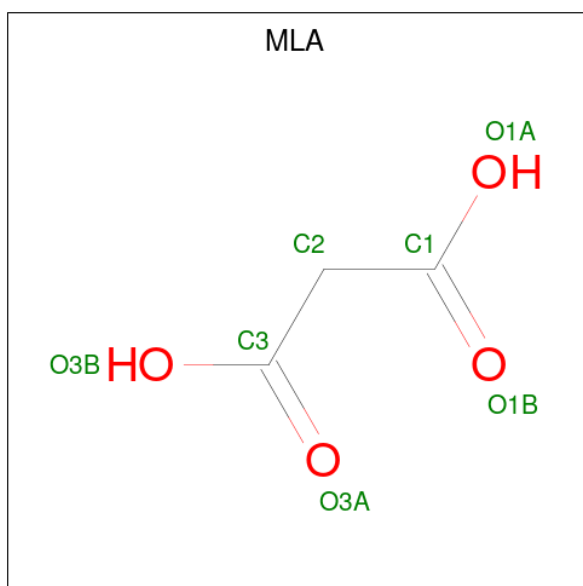
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	Fe	N	O	
			43	34	1	4	4	
							0	0

- Molecule 4 is 4-PHENYL-1H-IMIDAZOLE (CCD ID: PIM) (formula: $C_9H_8N_2$).



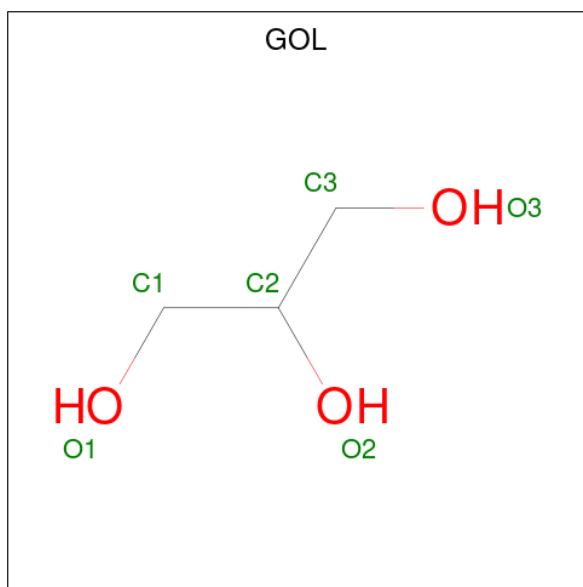
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	N		
			11	9	2		
						0	0

- Molecule 5 is MALONIC ACID (CCD ID: MLA) (formula: $C_3H_4O_4$).



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

- Molecule 6 is GLYCEROL (CCD ID: GOL) (formula: $C_3H_8O_3$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	A	1	Total	C	O	0	0
			6	3	3		

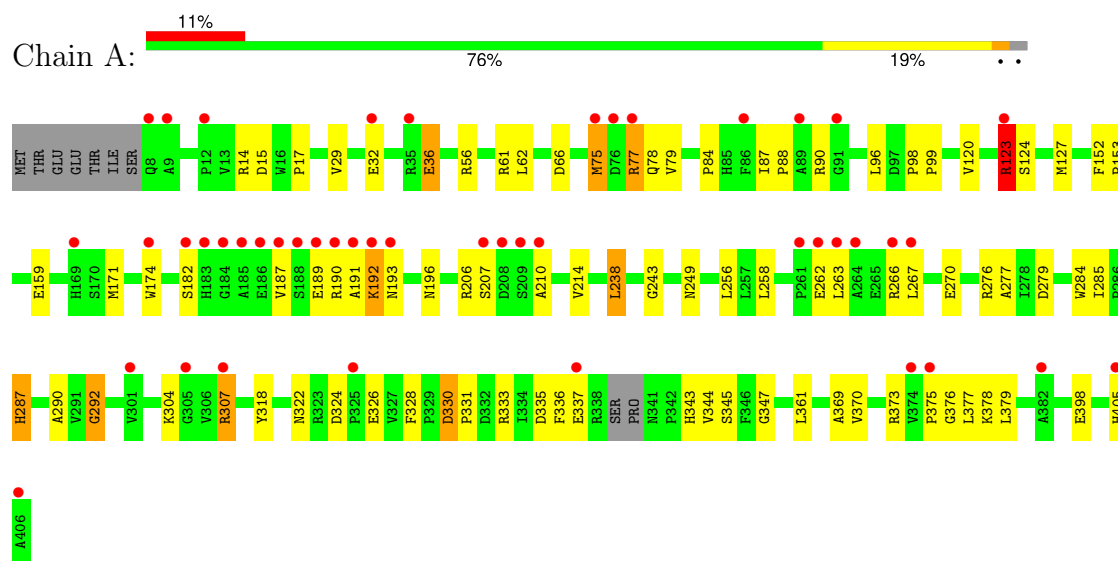
- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	326	Total	O	0	0
			326	326		

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: putative cytochrome P450



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	54.25Å 65.55Å 104.25Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	29.26 – 1.50 29.26 – 1.50	Depositor EDS
% Data completeness (in resolution range)	89.7 (29.26-1.50) 89.8 (29.26-1.50)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	0.06	Depositor
$\langle I/\sigma(I) \rangle$ ¹	0.88 (at 1.50Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.202 , 0.216 0.201 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å ²)	17.5	Xtriage
Anisotropy	0.137	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 34.7	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	0.96	EDS
Total number of atoms	3495	wwPDB-VP
Average B, all atoms (Å ²)	22.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.25% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: PIM, GOL, MLA, HEM, HG

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	0.41	1/3164 (0.0%)	0.91	15/4309 (0.3%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	99	PRO	CB-CG	-5.35	1.30	1.51

All (15) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	292	GLY	N-CA-C	11.70	120.85	110.21
1	A	287[A]	HIS	N-CA-C	7.64	120.50	111.71
1	A	287[B]	HIS	N-CA-C	7.64	120.50	111.71
1	A	98	PRO	N-CA-C	-7.16	103.59	110.47
1	A	77	ARG	N-CA-C	7.11	121.26	112.59
1	A	214	VAL	N-CA-C	6.34	116.51	110.42
1	A	284	TRP	N-CA-C	6.02	117.92	111.36
1	A	290	ALA	CB-CA-C	-5.48	110.24	116.54
1	A	330	ASP	CA-C-N	5.43	125.78	119.47
1	A	330	ASP	C-N-CA	5.43	125.78	119.47
1	A	398	GLU	N-CA-C	-5.20	105.68	112.23
1	A	75	MET	N-CA-C	5.18	117.67	111.71

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	66	ASP	CA-C-N	5.14	124.80	119.56
1	A	66	ASP	C-N-CA	5.14	124.80	119.56
1	A	36	GLU	N-CA-C	5.02	118.66	112.54

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	123	ARG	Sidechain

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	3093	0	3066	107	0
2	A	2	0	0	0	0
3	A	43	0	30	2	0
4	A	11	0	8	0	0
5	A	14	0	6	1	0
6	A	6	0	3	1	0
7	A	326	0	0	17	0
All	All	3495	0	3113	107	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 17.

All (107) 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:287[A]:HIS:HE1	7:A:523:HOH:O	1.01	1.33
1:A:61[B]:ARG:NH1	7:A:526:HOH:O	1.65	1.29
1:A:304:LYS:HE2	7:A:587:HOH:O	1.30	1.26
1:A:123:ARG:CD	1:A:159:GLU:HG2	1.72	1.19
1:A:287[B]:HIS:CE1	7:A:764:HOH:O	1.95	1.14
1:A:123:ARG:HD3	1:A:159:GLU:CD	1.76	1.11

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:287[B]:HIS:ND1	7:A:764:HOH:O	1.78	1.09
1:A:123:ARG:HD2	1:A:159:GLU:HG2	1.31	1.08
1:A:287[B]:HIS:H	1:A:287[B]:HIS:CD2	1.63	1.03
1:A:123:ARG:HD3	1:A:159:GLU:CG	1.90	1.02
1:A:123:ARG:CD	1:A:159:GLU:CG	2.42	0.97
1:A:123:ARG:HD3	1:A:159:GLU:HG2	1.45	0.94
1:A:287[B]:HIS:H	1:A:287[B]:HIS:HD2	1.11	0.94
1:A:322:ASN:HD21	1:A:344:VAL:H	1.15	0.89
1:A:32:GLU:OE2	7:A:531:HOH:O	1.94	0.85
1:A:56:ARG:NH2	1:A:326:GLU:HG3	1.92	0.84
1:A:287[B]:HIS:CD2	1:A:287[B]:HIS:N	2.45	0.83
1:A:56:ARG:HH21	1:A:326:GLU:HG3	1.43	0.83
1:A:189:GLU:HA	1:A:192:LYS:NZ	1.94	0.82
1:A:307:ARG:HH11	1:A:307:ARG:HB2	1.45	0.80
1:A:287[A]:HIS:CE1	7:A:523:HOH:O	1.88	0.80
1:A:123:ARG:HD3	1:A:159:GLU:OE1	1.82	0.79
1:A:78:GLN:OE1	7:A:594:HOH:O	1.98	0.79
1:A:192:LYS:NZ	1:A:192:LYS:HB2	1.98	0.79
1:A:266:ARG:HD2	1:A:270:GLU:OE1	1.83	0.78
1:A:190:ARG:HA	1:A:193:ASN:HD21	1.49	0.78
1:A:56:ARG:NH2	1:A:326:GLU:CG	2.49	0.76
1:A:87:ILE:HB	1:A:88:PRO:HD3	1.68	0.75
1:A:123:ARG:CD	1:A:159:GLU:CD	2.61	0.72
1:A:285:ILE:O	1:A:287[B]:HIS:CD2	2.43	0.71
1:A:190:ARG:HA	1:A:193:ASN:ND2	2.06	0.70
1:A:15[B]:ASP:OD1	7:A:622:HOH:O	2.09	0.70
1:A:343:HIS:HD2	1:A:345:SER:H	1.37	0.70
1:A:123:ARG:NH2	1:A:159:GLU:OE1	2.25	0.69
1:A:192:LYS:HB2	1:A:192:LYS:HZ3	1.56	0.69
1:A:285:ILE:O	1:A:287[B]:HIS:HD2	1.73	0.69
1:A:123:ARG:CD	1:A:159:GLU:OE1	2.40	0.69
1:A:328:PHE:O	1:A:331:PRO:HG3	1.95	0.67
1:A:123:ARG:HD2	1:A:159:GLU:CG	2.14	0.65
1:A:262:GLU:HG2	1:A:263:LEU:N	2.12	0.65
1:A:123:ARG:NE	1:A:159:GLU:OE1	2.29	0.65
1:A:262:GLU:HG2	1:A:263:LEU:HD12	1.78	0.65
1:A:187:VAL:O	1:A:192:LYS:HE3	1.97	0.64
1:A:189:GLU:HA	1:A:192:LYS:HZ1	1.62	0.63
1:A:56:ARG:HH22	1:A:326:GLU:HG2	1.63	0.63
1:A:243:GLY:HA2	3:A:430:HEM:HMC3	1.79	0.63
1:A:249:ASN:HD22	1:A:287[A]:HIS:HD2	1.48	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:120:VAL:HA	1:A:123:ARG:HH11	1.64	0.62
1:A:56:ARG:NH2	1:A:324:ASP:OD1	2.33	0.61
1:A:187:VAL:HG12	1:A:191:ALA:HB3	1.84	0.60
1:A:171:MET:HE2	1:A:174:TRP:HZ3	1.66	0.59
1:A:14:ARG:HE	1:A:29:VAL:HG13	1.70	0.57
1:A:249:ASN:HD22	1:A:287[A]:HIS:CD2	2.22	0.57
1:A:262:GLU:HG2	1:A:263:LEU:H	1.70	0.56
1:A:123:ARG:CZ	1:A:159:GLU:OE1	2.53	0.56
1:A:56:ARG:NH1	7:A:757:HOH:O	2.38	0.56
1:A:333:ARG:HH11	1:A:333:ARG:HG2	1.70	0.56
1:A:304:LYS:CE	7:A:587:HOH:O	2.09	0.56
1:A:152:PHE:HB3	1:A:153:PRO:HD3	1.86	0.55
1:A:62:LEU:C	1:A:62:LEU:HD13	2.33	0.53
1:A:78:GLN:HG3	7:A:692:HOH:O	2.06	0.53
1:A:307:ARG:HB2	1:A:307:ARG:NH1	2.22	0.52
1:A:330:ASP:N	1:A:331:PRO:HD3	2.23	0.52
1:A:267:LEU:HD21	1:A:277:ALA:CB	2.39	0.52
1:A:256:LEU:C	1:A:256:LEU:HD13	2.35	0.51
1:A:75:MET:HE1	1:A:90:ARG:HB2	1.93	0.51
1:A:120:VAL:HG12	1:A:123:ARG:NH1	2.24	0.51
1:A:279:ASP:OD2	1:A:343:HIS:HE1	1.94	0.51
1:A:376:GLY:HA3	1:A:405:HIS:CD2	2.46	0.51
1:A:84:PRO:HD2	1:A:182:SER:O	2.11	0.51
1:A:77:ARG:HG3	1:A:79:VAL:HG23	1.94	0.50
1:A:171:MET:HE2	1:A:174:TRP:CZ3	2.45	0.50
1:A:75:MET:SD	1:A:90:ARG:HA	2.52	0.49
1:A:361:LEU:C	1:A:361:LEU:HD23	2.37	0.49
1:A:14:ARG:HH11	1:A:14:ARG:HG2	1.76	0.49
1:A:262:GLU:HG2	1:A:263:LEU:CD1	2.42	0.49
1:A:238:LEU:CD2	7:A:754:HOH:O	2.60	0.48
1:A:276:ARG:HD2	1:A:336:PHE:O	2.12	0.48
1:A:333:ARG:NH1	1:A:335:ASP:HB2	2.29	0.48
1:A:318:TYR:CZ	1:A:347:GLY:HA2	2.49	0.48
1:A:189:GLU:HA	1:A:192:LYS:HZ2	1.76	0.48
1:A:287[B]:HIS:HD2	1:A:287[B]:HIS:N	1.92	0.47
1:A:206:ARG:O	1:A:207:SER:C	2.57	0.47
1:A:243:GLY:HA2	3:A:430:HEM:CMC	2.45	0.47
1:A:123:ARG:HG2	1:A:124:SER:N	2.29	0.46
1:A:210:ALA:HA	7:A:563:HOH:O	2.14	0.46
1:A:276:ARG:NH2	1:A:337:GLU:HG3	2.31	0.46
1:A:373:ARG:O	1:A:375:PRO:HD3	2.15	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:75:MET:HG3	7:A:741:HOH:O	2.16	0.45
1:A:287[B]:HIS:HE1	1:A:345:SER:HA	1.82	0.45
1:A:287[B]:HIS:CD2	7:A:469:HOH:O	2.69	0.45
1:A:369:ALA:HA	6:A:434:GOL:H12	1.99	0.45
1:A:78:GLN:N	1:A:78:GLN:CD	2.75	0.45
1:A:322:ASN:ND2	1:A:344:VAL:H	1.97	0.44
1:A:238:LEU:HD21	7:A:637:HOH:O	2.18	0.44
1:A:292:GLY:H	5:A:433:MLA:C1	2.32	0.43
1:A:238:LEU:HD13	1:A:238:LEU:HA	1.78	0.43
1:A:15[B]:ASP:OD2	1:A:17:PRO:HD3	2.18	0.43
1:A:32:GLU:O	1:A:36:GLU:HG3	2.18	0.43
1:A:370:VAL:CG1	1:A:377:LEU:HD22	2.49	0.43
1:A:77:ARG:H	1:A:77:ARG:HG2	1.53	0.42
1:A:258:LEU:HB3	1:A:379:LEU:HG	2.01	0.42
1:A:123:ARG:O	1:A:127:MET:HG3	2.19	0.41
1:A:189:GLU:HA	1:A:189:GLU:OE1	2.21	0.41
1:A:333:ARG:HG2	1:A:333:ARG:NH1	2.35	0.41
1:A:333:ARG:HH12	1:A:335:ASP:HA	1.86	0.40
1:A:377:LEU:HG	1:A:378:LYS:N	2.35	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	397/406 (98%)	386 (97%)	11 (3%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains [i](#)

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
1	A	326/335 (97%)	320 (98%)	6 (2%)	51 24

All (6) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	96	LEU
1	A	123	ARG
1	A	192	LYS
1	A	196	ASN
1	A	238	LEU
1	A	307	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	78	GLN
1	A	193	ASN
1	A	249	ASN
1	A	322	ASN
1	A	343	HIS
1	A	405	HIS

5.3.3 RNA [i](#)

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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 7 ligands modelled in this entry, 2 are monoatomic - leaving 5 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
4	PIM	A	431	3	11,12,12	2.49	6 (54%)	14,15,15	1.50	2 (14%)
3	HEM	A	430	4,1	50,50,50	1.63	9 (18%)	67,82,82	0.97	3 (4%)
6	GOL	A	434	-	5,5,5	5.08	5 (100%)	5,5,5	6.14	3 (60%)
5	MLA	A	433	2	6,6,6	4.32	3 (50%)	7,7,7	2.22	3 (42%)
5	MLA	A	432	-	6,6,6	4.18	3 (50%)	7,7,7	1.97	2 (28%)

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	PIM	A	431	3	-	0/4/4/4	0/2/2/2
3	HEM	A	430	4,1	-	4/14/54/54	-
6	GOL	A	434	-	-	2/4/4/4	-
5	MLA	A	433	2	-	2/4/4/4	-
5	MLA	A	432	-	-	0/4/4/4	-

All (26) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	A	433	MLA	O1A-C1	9.11	1.60	1.30
5	A	432	MLA	O1A-C1	8.82	1.60	1.30
6	A	434	GOL	C3-C2	-7.84	1.21	1.51
6	A	434	GOL	O3-C3	-5.40	1.19	1.42

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	430	HEM	CBB-CAB	4.92	1.54	1.30
3	A	430	HEM	CBC-CAC	4.66	1.52	1.30
6	A	434	GOL	O1-C1	4.52	1.61	1.42
5	A	432	MLA	O1B-C1	4.14	1.35	1.22
4	A	431	PIM	C8-C7	3.57	1.45	1.38
5	A	433	MLA	C2-C1	3.56	1.56	1.51
3	A	430	HEM	CAB-C3B	3.27	1.56	1.47
5	A	433	MLA	O1B-C1	3.26	1.32	1.22
4	A	431	PIM	C7-C6	3.20	1.44	1.39
4	A	431	PIM	C5-N1	-3.10	1.34	1.38
6	A	434	GOL	C1-C2	-3.06	1.40	1.51
4	A	431	PIM	C6-C5	2.96	1.52	1.47
6	A	434	GOL	O2-C2	-2.93	1.34	1.43
3	A	430	HEM	FE-NA	2.71	2.04	1.95
3	A	430	HEM	FE-NB	2.70	2.03	1.94
4	A	431	PIM	C10-C11	2.55	1.43	1.38
3	A	430	HEM	CAC-C3C	2.50	1.54	1.47
3	A	430	HEM	CMC-C2C	2.47	1.55	1.50
5	A	432	MLA	C2-C1	2.45	1.54	1.51
4	A	431	PIM	C4-N3	-2.40	1.33	1.37
3	A	430	HEM	CMD-C2D	2.19	1.55	1.50
3	A	430	HEM	O2A-CGA	-2.17	1.23	1.30

All (13) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	A	434	GOL	O3-C3-C2	11.24	160.99	110.38
6	A	434	GOL	O2-C2-C3	7.15	138.79	109.18
5	A	433	MLA	O1A-C1-C2	3.94	126.71	114.51
5	A	432	MLA	O1A-C1-C2	3.29	124.72	114.51
6	A	434	GOL	O1-C1-C2	3.29	125.18	110.38
5	A	433	MLA	O1B-C1-C2	-3.27	112.80	122.11
5	A	432	MLA	O1B-C1-C2	-3.12	113.23	122.11
3	A	430	HEM	CBA-CAA-C2A	2.62	119.77	112.53
3	A	430	HEM	CBC-CAC-C3C	-2.57	114.70	127.53
5	A	433	MLA	C3-C2-C1	2.38	121.36	112.95
3	A	430	HEM	O1A-CGA-CBA	-2.29	115.84	123.09
4	A	431	PIM	C4-N3-C2	2.16	110.18	105.94
4	A	431	PIM	C9-C10-C11	-2.15	117.58	120.24

There are no chirality outliers.

All (8) torsion outliers are listed below:

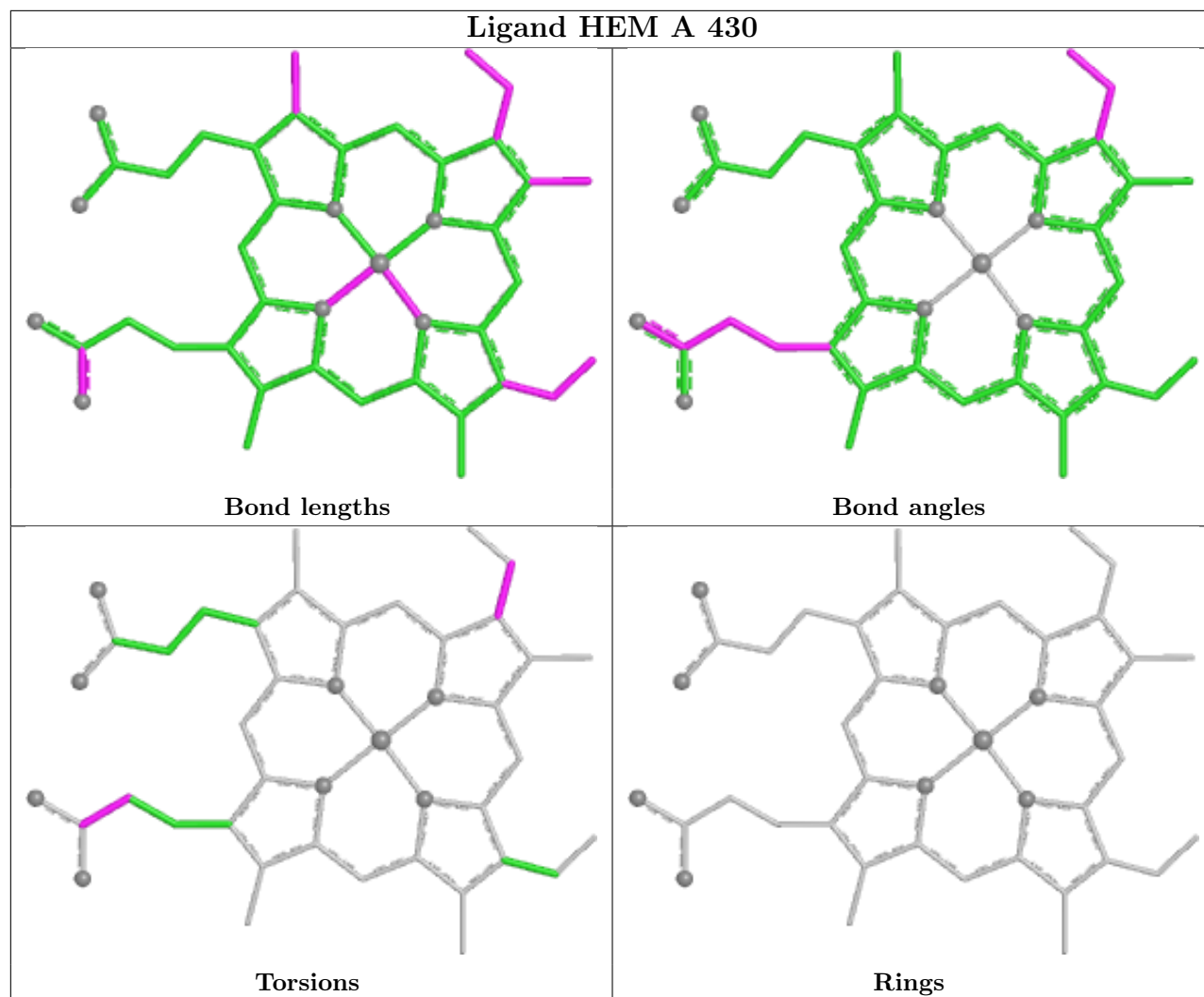
Mol	Chain	Res	Type	Atoms
3	A	430	HEM	C2C-C3C-CAC-CBC
3	A	430	HEM	C4C-C3C-CAC-CBC
6	A	434	GOL	C1-C2-C3-O3
6	A	434	GOL	O1-C1-C2-O2
3	A	430	HEM	CAA-CBA-CGA-O1A
3	A	430	HEM	CAA-CBA-CGA-O2A
5	A	433	MLA	C1-C2-C3-O3A
5	A	433	MLA	C1-C2-C3-O3B

There are no ring outliers.

3 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	430	HEM	2	0
6	A	434	GOL	1	0
5	A	433	MLA	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.



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	397/406 (97%)	0.71	46 (11%) 9 9	9, 20, 40, 52	4 (1%)

All (46) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	406	ALA	5.4
1	A	86	PHE	4.7
1	A	208	ASP	4.5
1	A	263	LEU	4.5
1	A	185	ALA	4.1
1	A	190	ARG	4.0
1	A	210	ALA	3.9
1	A	209	SER	3.8
1	A	187	VAL	3.7
1	A	8	GLN	3.7
1	A	184	GLY	3.6
1	A	192	LYS	3.6
1	A	261	PRO	3.2
1	A	264	ALA	3.1
1	A	207	SER	3.0
1	A	301	VAL	3.0
1	A	191	ALA	3.0
1	A	77	ARG	3.0
1	A	305	GLY	3.0
1	A	183	HIS	3.0
1	A	186	GLU	2.9
1	A	375	PRO	2.8
1	A	262	GLU	2.8
1	A	123	ARG	2.7
1	A	188	SER	2.7
1	A	193	ASN	2.6
1	A	189	GLU	2.6

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Mol	Chain	Res	Type	RSRZ
1	A	266	ARG	2.6
1	A	12	PRO	2.6
1	A	174	TRP	2.6
1	A	35	ARG	2.6
1	A	75	MET	2.5
1	A	89	ALA	2.4
1	A	76	ASP	2.4
1	A	91	GLY	2.4
1	A	267	LEU	2.4
1	A	182	SER	2.3
1	A	337	GLU	2.3
1	A	9	ALA	2.2
1	A	32	GLU	2.2
1	A	307	ARG	2.2
1	A	169	HIS	2.2
1	A	405	HIS	2.2
1	A	374	VAL	2.1
1	A	325	PRO	2.0
1	A	382	ALA	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 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.

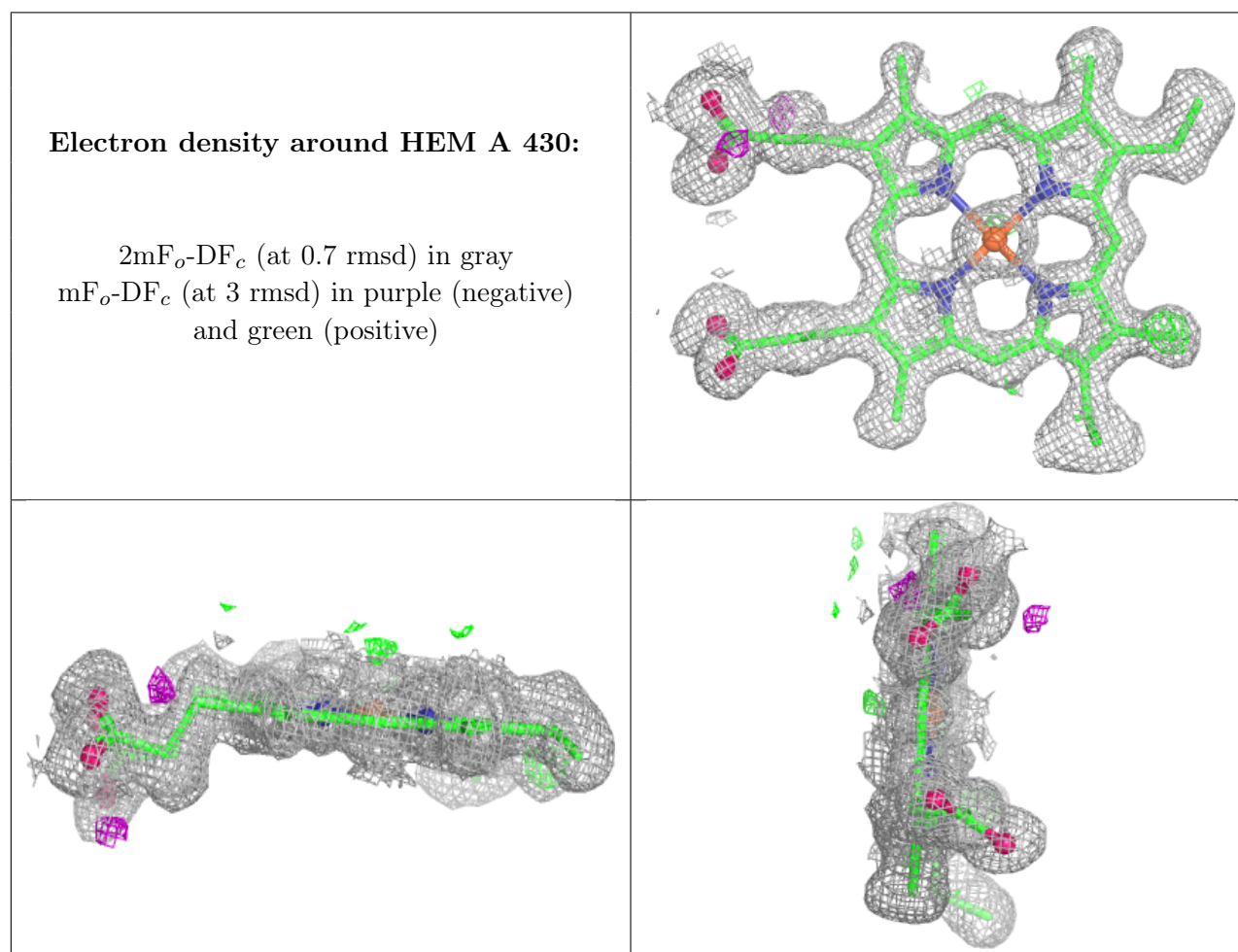
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
5	MLA	A	433	7/7	0.73	0.14	34,36,38,38	0
6	GOL	A	434	6/6	0.80	0.11	29,36,37,38	0
4	PIM	A	431	11/11	0.91	0.10	15,17,25,26	0
5	MLA	A	432	7/7	0.93	0.08	19,23,27,28	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	HEM	A	430	43/43	0.97	0.08	10,12,15,20	0
2	HG	A	436	1/1	0.99	0.07	17,17,17,17	1
2	HG	A	435	1/1	1.00	0.12	14,14,14,14	1

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.