



Full wwPDB X-ray Structure Validation Report ⓘ

Apr 5, 2026 – 12:00 PM UTC

PDB ID : 1B8F / pdb_00001b8f
Title : Histidine ammonia-lyase (HAL) from *Pseudomonas putida*
Authors : Schwede, T.F.; Schulz, G.E.
Deposited on : 1999-01-31
Resolution : 2.10 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
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

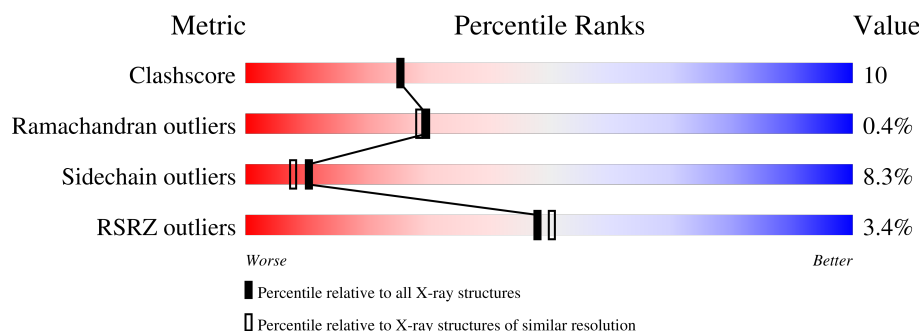
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.10 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	7164 (2.10-2.10)
Ramachandran outliers	187476	7099 (2.10-2.10)
Sidechain outliers	187428	7100 (2.10-2.10)
RSRZ outliers	180081	6662 (2.10-2.10)

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	507	3% 75% 20% ..

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
3	GOL	A	511	-	-	X	-

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 3983 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.

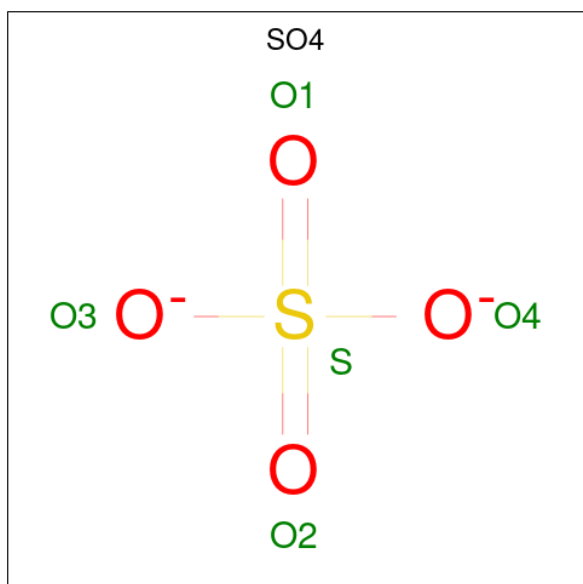
- Molecule 1 is a protein called Histidine ammonia-lyase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	507	3762	2360	670	716	16	0	0	0

There are 4 discrepancies between the modelled and reference sequences:

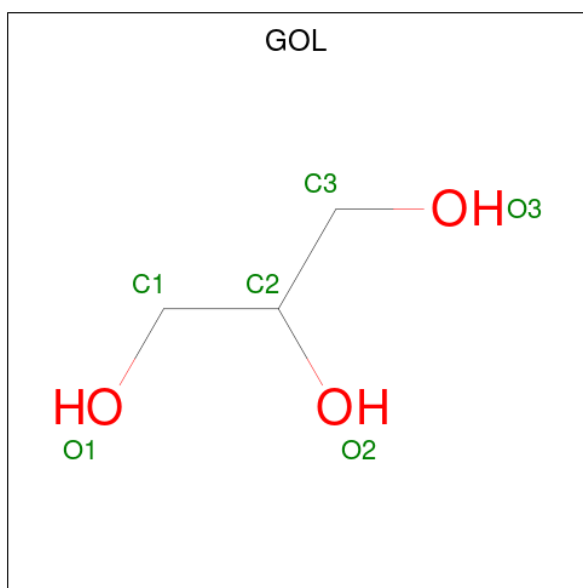
Chain	Residue	Modelled	Actual	Comment	Reference
A	142	MDO	ALA	chromophore	UNP P21310
A	142	MDO	SER	chromophore	UNP P21310
A	142	MDO	GLY	chromophore	UNP P21310
A	273	ALA	CYS	engineered mutation	UNP P21310

- Molecule 2 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



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

- Molecule 3 is GLYCEROL (CCD ID: GOL) (formula: $C_3H_8O_3$).



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

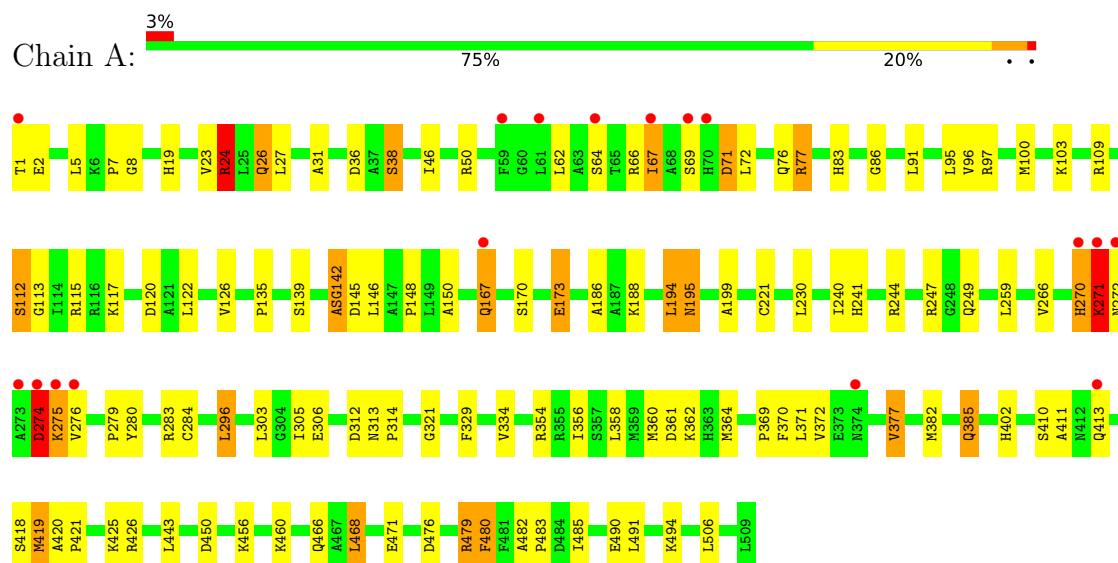
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	210	Total	O	0	0
			210	210		

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: Histidine ammonia-lyase



4 Data and refinement statistics

Property	Value	Source
Space group	I 2 2 2	Depositor
Cell constants a, b, c, α , β , γ	79.27Å 116.79Å 129.53Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	39.00 – 2.10 39.00 – 2.10	Depositor EDS
% Data completeness (in resolution range)	83.0 (39.00-2.10) 81.7 (39.00-2.10)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	8.20	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.48 (at 1.87Å)	Xtriage
Refinement program	REFMAC	Depositor
R, R_{free}	0.187 , 0.263 0.194 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å ²)	14.9	Xtriage
Anisotropy	1.685	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 45.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.45$, $\langle L^2 \rangle = 0.28$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	3983	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 5.18% 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: SO4, GOL, MDO

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	1.06	0/3806	1.62	33/5161 (0.6%)

There are no bond length outliers.

All (33) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	270	HIS	CA-CB-CG	12.91	126.71	113.80
1	A	313	ASN	CA-C-O	-10.73	109.35	119.76
1	A	450	ASP	CA-CB-CG	9.16	121.76	112.60
1	A	479	ARG	CD-NE-CZ	8.72	136.60	124.40
1	A	24	ARG	CD-NE-CZ	7.42	134.78	124.40
1	A	240	ILE	CB-CA-C	-7.29	102.64	111.97
1	A	303	LEU	CA-C-N	6.94	127.63	120.00
1	A	303	LEU	C-N-CA	6.94	127.63	120.00
1	A	321	GLY	CA-C-N	6.85	131.21	121.42
1	A	321	GLY	C-N-CA	6.85	131.21	121.42
1	A	480	PHE	CA-CB-CG	6.79	120.59	113.80
1	A	36	ASP	CA-CB-CG	6.75	119.35	112.60
1	A	402	HIS	N-CA-C	-6.66	105.31	113.50
1	A	5	LEU	N-CA-C	6.40	119.50	109.96
1	A	77	ARG	CD-NE-CZ	-6.33	115.54	124.40
1	A	377	VAL	CB-CA-C	6.06	117.16	110.62
1	A	120	ASP	CA-CB-CG	5.94	118.54	112.60
1	A	274	ASP	CA-CB-CG	5.85	118.45	112.60
1	A	485	ILE	O-C-N	5.84	127.98	121.90
1	A	221	CYS	CA-C-N	5.76	126.37	119.98
1	A	221	CYS	C-N-CA	5.76	126.37	119.98
1	A	115	ARG	CD-NE-CZ	5.72	132.41	124.40
1	A	377	VAL	N-CA-CB	-5.69	105.56	112.33
1	A	19	HIS	CA-CB-CG	5.58	119.39	113.80

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	71	ASP	CA-CB-CG	5.56	118.16	112.60
1	A	476	ASP	CA-C-N	5.42	131.13	121.75
1	A	476	ASP	C-N-CA	5.42	131.13	121.75
1	A	241	HIS	CA-CB-CG	-5.42	108.38	113.80
1	A	354	ARG	NE-CZ-NH2	5.42	124.07	119.20
1	A	314	PRO	N-CA-CB	5.38	108.52	102.60
1	A	283	ARG	CD-NE-CZ	5.28	131.79	124.40
1	A	77	ARG	NE-CZ-NH1	-5.12	116.38	121.50
1	A	83	HIS	CA-CB-CG	-5.12	108.68	113.80

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3762	0	3814	73	0
2	A	5	0	0	0	0
3	A	6	0	8	6	0
4	A	210	0	0	9	0
All	All	3983	0	3822	73	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 10.

All (73) 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:482:ALA:HB3	1:A:483:PRO:HD3	1.61	0.79
1:A:471:GLU:HG3	1:A:491:LEU:HD11	1.66	0.76
1:A:271:LYS:HG2	1:A:272:ASN:H	1.50	0.76
1:A:360:MET:HE3	1:A:370:PHE:HA	1.73	0.70
1:A:76:GLN:HE22	3:A:511:GOL:H31	1.56	0.69
1:A:150:ALA:HB2	1:A:194:LEU:CD1	2.25	0.67
1:A:490:GLU:OE2	1:A:494:LYS:HE2	1.98	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:145:ASP:HB3	1:A:148:PRO:HG2	1.80	0.62
1:A:356:ILE:HG22	1:A:385:GLN:HG3	1.83	0.61
1:A:356:ILE:O	1:A:360:MET:HG2	2.00	0.60
1:A:139:SER:HB3	1:A:421:PRO:HG2	1.82	0.59
1:A:72:LEU:HD11	3:A:511:GOL:H32	1.84	0.59
1:A:76:GLN:HE22	3:A:511:GOL:C3	2.15	0.58
1:A:259:LEU:CD2	1:A:460:LYS:HD2	2.33	0.58
1:A:271:LYS:CG	1:A:272:ASN:H	2.16	0.56
1:A:76:GLN:NE2	3:A:511:GOL:H31	2.21	0.55
1:A:188:LYS:H	3:A:511:GOL:C3	2.19	0.55
1:A:443:LEU:HD23	1:A:468:LEU:HD11	1.89	0.55
1:A:506:LEU:HD12	4:A:689:HOH:O	2.05	0.55
1:A:112:SER:HB3	1:A:312:ASP:OD2	2.08	0.54
1:A:91:LEU:HD22	1:A:95:LEU:HD23	1.90	0.54
1:A:96:VAL:O	1:A:100:MET:HG3	2.08	0.54
1:A:370:PHE:CE1	1:A:382:MET:HG3	2.44	0.53
1:A:139:SER:OG	1:A:145:ASP:HA	2.09	0.52
1:A:296:LEU:C	1:A:296:LEU:HD12	2.35	0.52
1:A:150:ALA:HB2	1:A:194:LEU:HD11	1.92	0.52
1:A:259:LEU:HD23	1:A:460:LYS:HD2	1.91	0.52
1:A:275:LYS:HE2	4:A:612:HOH:O	2.10	0.52
1:A:26:GLN:HG2	4:A:626:HOH:O	2.11	0.51
1:A:360:MET:CE	1:A:370:PHE:HA	2.41	0.50
1:A:142:MDO:N3	1:A:329:PHE:HB2	2.25	0.50
1:A:186:ALA:HA	4:A:639:HOH:O	2.11	0.50
1:A:360:MET:CE	1:A:370:PHE:HD1	2.25	0.50
1:A:279:PRO:HG2	1:A:358:LEU:CD1	2.42	0.49
1:A:150:ALA:HB2	1:A:194:LEU:HD13	1.94	0.49
1:A:418:SER:O	1:A:419:MET:HB2	2.12	0.49
1:A:38:SER:OG	1:A:113:GLY:HA2	2.12	0.49
1:A:274:ASP:O	1:A:275:LYS:C	2.56	0.49
1:A:170:SER:OG	1:A:173:GLU:HG3	2.13	0.48
1:A:479:ARG:HD3	4:A:596:HOH:O	2.14	0.47
1:A:188:LYS:N	3:A:511:GOL:O3	2.46	0.47
1:A:266:VAL:HG23	1:A:456:LYS:O	2.14	0.47
1:A:279:PRO:HG2	1:A:358:LEU:HD11	1.95	0.47
1:A:411:ALA:O	1:A:413:GLN:NE2	2.48	0.47
1:A:46:ILE:HA	4:A:575:HOH:O	2.15	0.47
1:A:86:GLY:HA3	1:A:135:PRO:HG2	1.98	0.46
1:A:230:LEU:HB3	1:A:270:HIS:CD2	2.51	0.46
1:A:420:ALA:HB3	1:A:421:PRO:HD3	1.98	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:362:LYS:HD3	1:A:369:PRO:HG3	1.99	0.45
1:A:142:MDO:CB2	1:A:146:LEU:HD11	2.47	0.45
1:A:145:ASP:CG	1:A:199:ALA:HB3	2.42	0.45
1:A:66:ARG:O	1:A:67:ILE:C	2.59	0.45
1:A:360:MET:HE3	1:A:370:PHE:HD1	1.81	0.44
1:A:103:LYS:NZ	1:A:195:ASN:O	2.50	0.44
1:A:270:HIS:O	1:A:271:LYS:C	2.60	0.44
1:A:167:GLN:HE21	1:A:167:GLN:HB3	1.52	0.43
1:A:280:TYR:O	1:A:284:CYS:HB2	2.18	0.43
1:A:456:LYS:HZ2	1:A:456:LYS:HG2	1.76	0.43
1:A:122:LEU:O	1:A:126:VAL:HG23	2.18	0.43
1:A:361:ASP:HA	4:A:567:HOH:O	2.19	0.42
1:A:306:GLU:HG2	1:A:334:VAL:CG2	2.49	0.42
1:A:8:GLY:HA3	1:A:109:ARG:HD2	2.01	0.42
1:A:77:ARG:HH11	1:A:77:ARG:HD3	1.59	0.41
1:A:23:VAL:HG22	1:A:24:ARG:N	2.35	0.41
1:A:480:PHE:CE1	1:A:482:ALA:HB3	2.55	0.41
1:A:244:ARG:HD2	1:A:249:GLN:OE1	2.21	0.41
1:A:360:MET:HE3	1:A:370:PHE:CD1	2.55	0.41
1:A:66:ARG:O	1:A:66:ARG:HG2	2.21	0.41
1:A:71:ASP:HB3	4:A:604:HOH:O	2.20	0.41
1:A:7:PRO:HG2	1:A:31:ALA:CB	2.50	0.41
1:A:50:ARG:HH11	1:A:50:ARG:HD2	1.66	0.41
1:A:1:THR:HA	4:A:671:HOH:O	2.20	0.40
1:A:23:VAL:HG12	1:A:97:ARG:HD2	2.03	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	502/507 (99%)	481 (96%)	19 (4%)	2 (0%)	30	28

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	271	LYS
1	A	275	LYS

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	387/387 (100%)	355 (92%)	32 (8%)	10 8

All (32) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	2	GLU
1	A	24	ARG
1	A	26	GLN
1	A	27	LEU
1	A	38	SER
1	A	62	LEU
1	A	64	SER
1	A	67	ILE
1	A	69	SER
1	A	112	SER
1	A	117	LYS
1	A	167	GLN
1	A	173	GLU
1	A	194	LEU
1	A	195	ASN
1	A	247	ARG
1	A	271	LYS
1	A	274	ASP
1	A	276	VAL
1	A	296	LEU
1	A	305	ILE
1	A	364	MET
1	A	371	LEU

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Mol	Chain	Res	Type
1	A	372	VAL
1	A	377	VAL
1	A	385	GLN
1	A	410	SER
1	A	419	MET
1	A	425	LYS
1	A	426	ARG
1	A	466	GLN
1	A	468	LEU

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	26	GLN
1	A	74	ASN
1	A	167	GLN
1	A	272	ASN
1	A	402	HIS
1	A	413	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

1 non-standard protein/DNA/RNA residue is 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	$\# Z > 2$
1	MDO	A	142	1	11,13,14	2.25	3 (27%)	15,18,20	7.31	9 (60%)

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
1	MDO	A	142	1	-	2/4/23/24	0/1/1/1

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	142	MDO	C1-N2	-4.43	1.25	1.32
1	A	142	MDO	CA2-N2	4.13	1.46	1.39
1	A	142	MDO	C2-N3	-2.08	1.35	1.40

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	142	MDO	C2-CA2-N2	-15.64	97.74	108.95
1	A	142	MDO	C3-CA3-N3	15.61	147.96	112.43
1	A	142	MDO	CA2-C2-N3	9.42	111.41	103.50
1	A	142	MDO	O2-C2-CA2	-7.74	126.08	131.02
1	A	142	MDO	CA2-N2-C1	7.69	112.28	105.39
1	A	142	MDO	CA3-N3-C2	-6.64	109.08	123.67
1	A	142	MDO	C1-CA1-N1	5.95	126.67	109.13
1	A	142	MDO	C2-N3-C1	-3.80	106.31	108.07
1	A	142	MDO	O3-C3-CA3	-2.21	115.50	125.47

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	A	142	MDO	C3-CA3-N3-C2
1	A	142	MDO	C3-CA3-N3-C1

There are no ring outliers.

1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	A	142	MDO	2	0

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

2 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	$\# Z > 2$
3	GOL	A	511	-	5,5,5	0.81	0	5,5,5	1.00	0
2	SO4	A	510	-	4,4,4	0.78	0	6,6,6	0.18	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
3	GOL	A	511	-	-	2/4/4/4	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	511	GOL	O1-C1-C2-C3
3	A	511	GOL	O1-C1-C2-O2

There are no ring outliers.

1 monomer is involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	511	GOL	6	0

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	506/507 (99%)	-0.13	17 (3%)	48 50	9, 19, 40, 65	0

All (17) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	1	THR	4.6
1	A	274	ASP	4.5
1	A	273	ALA	4.3
1	A	272	ASN	3.2
1	A	59	PHE	3.1
1	A	67	ILE	2.8
1	A	167	GLN	2.6
1	A	271	LYS	2.5
1	A	70	HIS	2.5
1	A	69	SER	2.4
1	A	374	ASN	2.4
1	A	64	SER	2.4
1	A	413	GLN	2.3
1	A	276	VAL	2.3
1	A	270	HIS	2.1
1	A	61	LEU	2.0
1	A	275	LYS	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [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
1	MDO	A	142	13/14	0.95	0.07	14,18,21,21	0

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
3	GOL	A	511	6/6	0.81	0.14	24,30,31,33	0
2	SO4	A	510	5/5	0.90	0.18	58,59,59,59	0

6.5 Other polymers [i](#)

There are no such residues in this entry.