



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 08:12 PM UTC

PDB ID : 9IRQ / pdb_00009irq
Title : Alginate lyase (Microbulbifer sp. ALW1)
Authors : Zhu, Y.B.
Deposited on : 2024-07-16
Resolution : 1.98 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
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

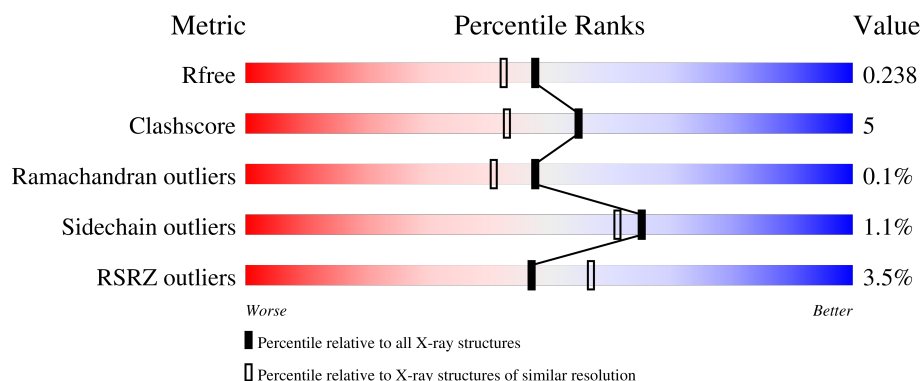
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.98 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	1506 (1.98-1.98)
Clashscore	190562	1534 (1.98-1.98)
Ramachandran outliers	187476	1518 (1.98-1.98)
Sidechain outliers	187428	1518 (1.98-1.98)
RSRZ outliers	180081	1506 (1.98-1.98)

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	707	4% 88% 12%
1	B	707	3% 89% 10%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 11464 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 Alginate lyase AlgL17.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	707	Total	C	N	O	S	0	0	0
			5538	3519	955	1057	7			
1	B	707	Total	C	N	O	S	0	0	0
			5548	3525	964	1052	7			

- Molecule 2 is ZINC ION (CCD ID: ZN) (formula: Zn) (labeled as "Ligand of Interest" by depositor).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	Zn	0	0
			1	1		
2	B	1	Total	Zn	0	0
			1	1		

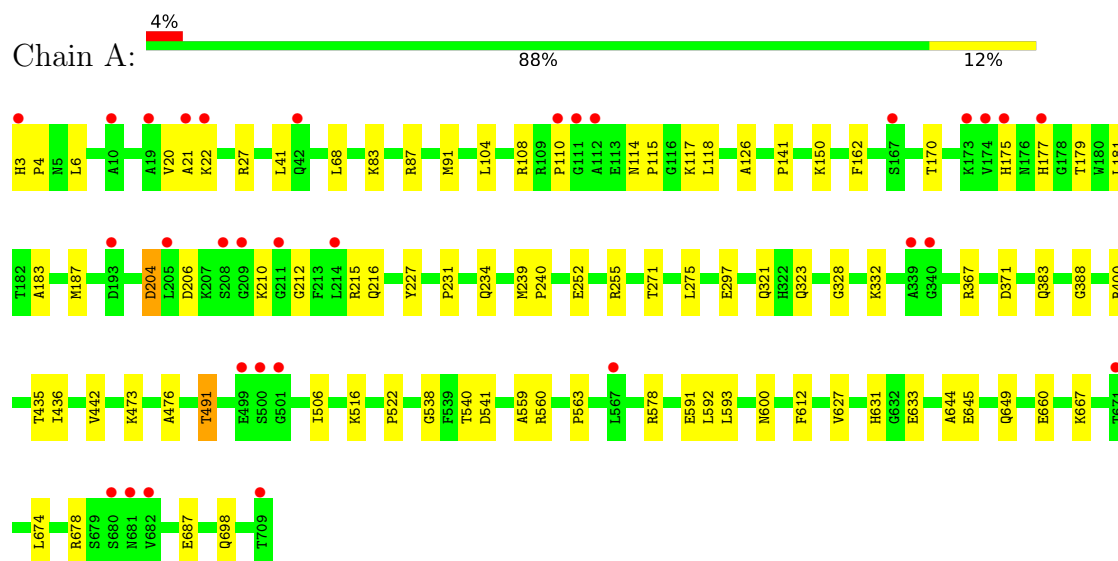
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	173	Total	O	0	0
			173	173		
3	B	203	Total	O	0	0
			203	203		

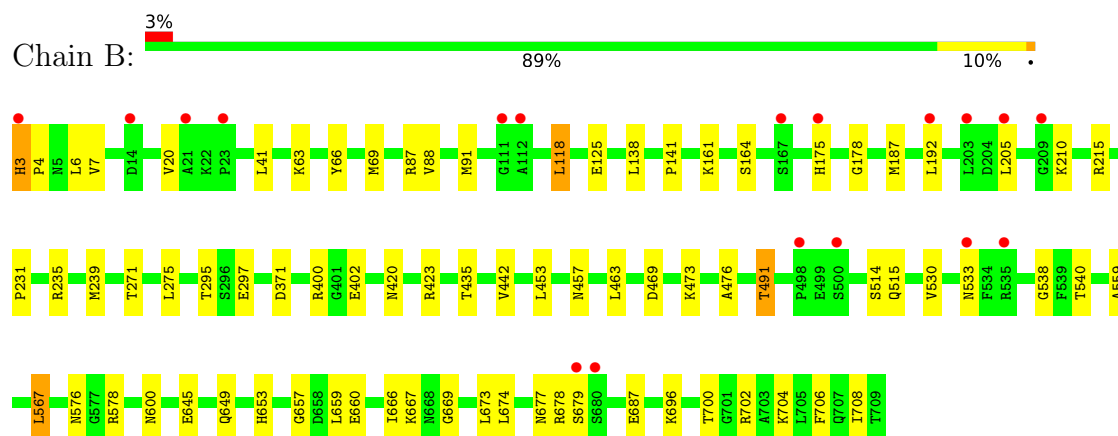
3 Residue-property plots [i](#)

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: Alginate lyase AlgL17



• Molecule 1: Alginate lyase AlgL17



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	87.11Å 90.42Å 88.29Å 90.00° 110.01° 90.00°	Depositor
Resolution (Å)	82.96 – 1.98 82.96 – 1.98	Depositor EDS
% Data completeness (in resolution range)	94.5 (82.96-1.98) 94.5 (82.96-1.98)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.27 (at 1.98Å)	Xtriage
Refinement program	PHENIX 1.12	Depositor
R, R_{free}	0.188 , 0.239 0.192 , 0.238	Depositor DCC
R_{free} test set	4006 reflections (4.50%)	wwPDB-VP
Wilson B-factor (Å ²)	28.2	Xtriage
Anisotropy	0.509	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 34.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.000 for l,-k,h	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	11464	wwPDB-VP
Average B, all atoms (Å ²)	32.0	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 60.47 % 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 1.5149e-05. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

¹ Intensities estimated from amplitudes.

² Theoretical 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: ZN

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	0.39	0/5671	0.61	1/7703 (0.0%)
1	B	0.63	0/5681	0.83	2/7713 (0.0%)
All	All	0.52	0/11352	0.73	3/15416 (0.0%)

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	3	HIS	CA-CB-CG	5.96	119.76	113.80
1	B	7	VAL	N-CA-C	-5.60	104.89	111.00
1	A	206	ASP	N-CA-C	-5.13	106.97	113.18

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	5538	0	5348	57	0
1	B	5548	0	5378	49	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	173	0	0	1	0
3	B	203	0	0	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
All	All	11464	0	10726	104	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 5.

All (104) 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:560:ARG:NE	1:A:591:GLU:OE1	1.95	0.98
1:A:649:GLN:HG3	1:A:667:LYS:HG2	1.50	0.91
1:B:87:ARG:HG3	1:B:91:MET:HE2	1.59	0.84
1:A:108:ARG:HD2	1:A:110:PRO:HG3	1.62	0.79
1:B:679:SER:O	1:B:702:ARG:NH1	2.14	0.77
1:A:108:ARG:HG3	1:A:115:PRO:HB3	1.67	0.77
1:A:20:VAL:O	1:A:27:ARG:NH1	2.17	0.77
1:B:3:HIS:HB2	1:B:141:PRO:HD3	1.71	0.72
1:A:126:ALA:HB2	1:A:181:LEU:HD23	1.73	0.71
1:B:533:ASN:ND2	1:B:567:LEU:O	2.24	0.70
1:A:87:ARG:HG3	1:A:91:MET:HE2	1.73	0.69
1:A:371:ASP:CG	1:A:578:ARG:HH22	2.01	0.68
1:A:114:ASN:OD1	1:A:170:THR:OG1	2.14	0.65
1:B:3:HIS:N	1:B:4:PRO:HD3	2.12	0.65
1:B:514:SER:OG	1:B:515:GLN:NE2	2.30	0.64
1:B:371:ASP:O	1:B:400:ARG:NH1	2.30	0.64
1:B:476:ALA:HB2	1:B:491:THR:HB	1.80	0.63
1:A:41:LEU:HD12	1:A:83:LYS:HE3	1.83	0.61
1:A:239:MET:HE3	1:A:297:GLU:HA	1.85	0.59
1:A:117:LYS:HE2	1:A:162:PHE:CD1	2.38	0.59
1:B:175:HIS:CD2	1:B:178:GLY:H	2.21	0.59
1:B:175:HIS:HD2	1:B:178:GLY:H	1.52	0.58
1:A:108:ARG:CG	1:A:115:PRO:HB3	2.33	0.57
1:A:600:ASN:HA	1:B:540:THR:OG1	2.05	0.57
1:A:476:ALA:HB2	1:A:491:THR:HB	1.87	0.56
1:A:506:ILE:HG12	1:A:627:VAL:HG22	1.88	0.56
1:B:453:LEU:CD2	1:B:457:ASN:ND2	2.69	0.56
1:B:469:ASP:O	1:B:473:LYS:HE3	2.07	0.54
1:B:371:ASP:CG	1:B:578:ARG:HH22	2.15	0.54
1:A:3:HIS:HB2	1:A:141:PRO:HD3	1.88	0.54
1:A:687:GLU:OE1	1:A:698:GLN:HB2	2.08	0.53
1:B:473:LYS:NZ	3:B:904:HOH:O	2.41	0.53
1:B:235:ARG:NH1	1:B:297:GLU:OE2	2.42	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:321:GLN:O	1:A:323:GLN:HG3	2.11	0.51
1:A:645:GLU:H	1:A:645:GLU:CD	2.19	0.51
1:B:3:HIS:N	1:B:4:PRO:CD	2.73	0.51
1:B:687:GLU:OE2	1:B:700:THR:HG22	2.11	0.50
1:B:659:LEU:HD23	1:B:677:ASN:HA	1.92	0.50
1:A:400:ARG:HH12	1:A:633:GLU:CD	2.19	0.50
1:A:538:GLY:HA2	1:A:559:ALA:HB2	1.93	0.50
1:B:6:LEU:HD13	1:B:187:MET:HE2	1.93	0.50
1:A:6:LEU:HD13	1:A:187:MET:HE2	1.93	0.49
1:B:420:ASN:OD1	1:B:423:ARG:NH1	2.44	0.49
1:A:108:ARG:HD3	1:A:115:PRO:HB3	1.94	0.49
1:B:649:GLN:HG3	1:B:667:LYS:HG2	1.94	0.49
1:A:68:LEU:HD12	1:A:91:MET:HE1	1.95	0.49
1:B:666:ILE:HG22	1:B:669:GLY:HA3	1.94	0.49
1:B:271:THR:O	1:B:275:LEU:HG	2.13	0.48
1:B:673:LEU:HD22	1:B:708:ILE:HD11	1.95	0.48
1:B:657:GLY:O	1:B:678:ARG:NH2	2.42	0.48
1:A:436:ILE:HD12	1:A:522:PRO:HG3	1.95	0.48
1:B:69:MET:HG2	1:B:91:MET:HE3	1.94	0.48
1:A:540:THR:OG1	1:B:600:ASN:HA	2.14	0.47
1:A:3:HIS:N	1:A:4:PRO:HD2	2.29	0.47
1:A:231:PRO:HG3	1:A:275:LEU:HD13	1.96	0.47
1:A:21:ALA:HB3	1:A:22:LYS:HE2	1.97	0.47
1:A:3:HIS:CB	1:A:141:PRO:HD3	2.46	0.46
1:A:179:THR:HG22	1:A:240:PRO:HB2	1.97	0.46
1:A:108:ARG:HE	1:A:108:ARG:HB3	1.62	0.45
1:B:235:ARG:HD3	1:B:295:THR:HG21	1.98	0.45
1:A:367:ARG:O	1:A:473:LYS:HE2	2.16	0.45
1:A:210:LYS:O	1:A:215:ARG:HG2	2.16	0.45
1:A:435:THR:HG22	1:A:442:VAL:HG22	1.98	0.45
1:B:63:LYS:HA	1:B:66:TYR:CE2	2.52	0.45
1:B:210:LYS:O	1:B:215:ARG:HG2	2.17	0.45
1:A:108:ARG:CD	1:A:110:PRO:HG3	2.42	0.44
1:B:3:HIS:CB	1:B:141:PRO:HD3	2.44	0.44
1:A:563:PRO:HG3	1:A:592:LEU:HD12	1.99	0.44
1:A:175:HIS:NE2	1:A:177:HIS:HB3	2.32	0.44
1:A:660:GLU:CG	1:A:678:ARG:HD2	2.47	0.44
1:A:227:TYR:HB3	1:A:234:GLN:HB2	1.99	0.44
1:A:593:LEU:O	1:A:612:PHE:HA	2.18	0.44
1:A:3:HIS:N	1:A:4:PRO:CD	2.81	0.43
1:B:4:PRO:HB3	1:B:138:LEU:HD23	2.00	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:161:LYS:HE2	1:B:161:LYS:HB3	1.60	0.43
1:B:538:GLY:HA2	1:B:559:ALA:HB2	2.00	0.43
1:B:118:LEU:HD12	1:B:118:LEU:HA	1.84	0.43
1:B:567:LEU:HD23	1:B:567:LEU:HA	1.91	0.43
1:B:704:LYS:HG2	1:B:706:PHE:CZ	2.53	0.43
1:A:332:LYS:HA	1:A:332:LYS:HD3	1.67	0.43
1:B:41:LEU:HD22	1:B:87:ARG:HG2	2.00	0.43
1:B:435:THR:HG22	1:B:442:VAL:HG22	2.00	0.43
1:A:383:GLN:O	1:A:388:GLY:HA3	2.19	0.43
1:A:104:LEU:HA	1:A:104:LEU:HD23	1.79	0.43
1:A:540:THR:HG23	1:A:541:ASP:OD1	2.19	0.42
1:A:212:GLY:O	1:A:216:GLN:HG3	2.19	0.42
1:B:88:VAL:HA	1:B:91:MET:HE3	2.02	0.42
1:A:183:ALA:O	1:A:187:MET:HG3	2.19	0.42
1:B:402:GLU:HG3	1:B:576:ASN:HB3	2.02	0.42
1:A:252:GLU:CD	1:A:255:ARG:HE	2.27	0.41
1:A:631:HIS:HE1	1:A:644:ALA:O	2.03	0.41
1:B:653:HIS:NE2	1:B:660:GLU:OE2	2.52	0.41
1:B:231:PRO:HG3	1:B:275:LEU:HD13	2.02	0.41
1:A:560:ARG:CD	1:A:591:GLU:OE1	2.67	0.41
1:B:164:SER:HB2	1:B:205:LEU:HD21	2.02	0.41
1:B:645:GLU:H	1:B:645:GLU:CD	2.28	0.41
1:A:108:ARG:CD	1:A:115:PRO:HB3	2.51	0.41
1:B:696:LYS:HE2	1:B:696:LYS:HB2	1.90	0.40
1:A:150:LYS:HD3	1:A:150:LYS:HA	1.88	0.40
1:A:516:LYS:HB3	1:A:516:LYS:HE3	1.93	0.40
1:B:453:LEU:HD23	1:B:453:LEU:O	2.21	0.40
1:A:271:THR:O	1:A:275:LEU:HG	2.22	0.40
1:A:328:GLY:N	3:A:918:HOH:O	2.55	0.40
1:B:239:MET:HE3	1:B:297:GLU:HG3	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	705/707 (100%)	685 (97%)	19 (3%)	1 (0%)	48	41
1	B	705/707 (100%)	681 (97%)	24 (3%)	0	100	100
All	All	1410/1414 (100%)	1366 (97%)	43 (3%)	1 (0%)	48	41

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	204	ASP

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	576/582 (99%)	572 (99%)	4 (1%)	76	73
1	B	577/582 (99%)	568 (98%)	9 (2%)	55	50
All	All	1153/1164 (99%)	1140 (99%)	13 (1%)	65	61

All (13) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	118	LEU
1	A	204	ASP
1	A	491	THR
1	A	674	LEU
1	B	20	VAL
1	B	118	LEU
1	B	125	GLU
1	B	192	LEU
1	B	463	LEU
1	B	491	THR
1	B	530	VAL
1	B	567	LEU

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Mol	Chain	Res	Type
1	B	674	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (13) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	42	GLN
1	A	61	GLN
1	A	121	GLN
1	A	459	ASN
1	A	518	GLN
1	A	549	ASN
1	A	672	GLN
1	B	61	GLN
1	B	121	GLN
1	B	153	GLN
1	B	175	HIS
1	B	218	ASN
1	B	515	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

Of 2 ligands modelled in this entry, 2 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

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	707/707 (100%)	0.43	31 (4%)	39	49	19, 31, 51, 64	0
1	B	707/707 (100%)	0.37	18 (2%)	58	68	20, 30, 47, 63	0
All	All	1414/1414 (100%)	0.40	49 (3%)	47	57	19, 31, 49, 64	0

All (49) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	21	ALA	3.3
1	B	3	HIS	3.3
1	A	175	HIS	3.2
1	A	709	THR	3.1
1	B	203	LEU	3.1
1	A	209	GLY	3.0
1	A	193	ASP	2.8
1	A	211	GLY	2.7
1	A	174	VAL	2.7
1	A	340	GLY	2.7
1	B	192	LEU	2.7
1	A	3	HIS	2.7
1	B	21	ALA	2.6
1	A	501	GLY	2.6
1	A	671	THR	2.6
1	B	533	ASN	2.6
1	A	110	PRO	2.6
1	B	209	GLY	2.6
1	A	22	LYS	2.5
1	A	681	ASN	2.4
1	A	167	SER	2.4
1	A	10	ALA	2.4
1	A	208	SER	2.4
1	A	111	GLY	2.4

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Mol	Chain	Res	Type	RSRZ
1	A	680	SER	2.4
1	A	177	HIS	2.3
1	B	14	ASP	2.3
1	B	500	SER	2.3
1	A	173	LYS	2.3
1	A	19	ALA	2.2
1	B	111	GLY	2.2
1	B	167	SER	2.2
1	B	535	ARG	2.2
1	B	498	PRO	2.2
1	A	567	LEU	2.2
1	A	112	ALA	2.2
1	B	112	ALA	2.2
1	B	205	LEU	2.1
1	A	499	GLU	2.1
1	A	339	ALA	2.1
1	B	680	SER	2.1
1	B	23	PRO	2.1
1	A	42	GLN	2.1
1	A	500	SER	2.1
1	B	679	SER	2.1
1	A	682	VAL	2.1
1	A	205	LEU	2.1
1	B	175	HIS	2.0
1	A	214	LEU	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](#)

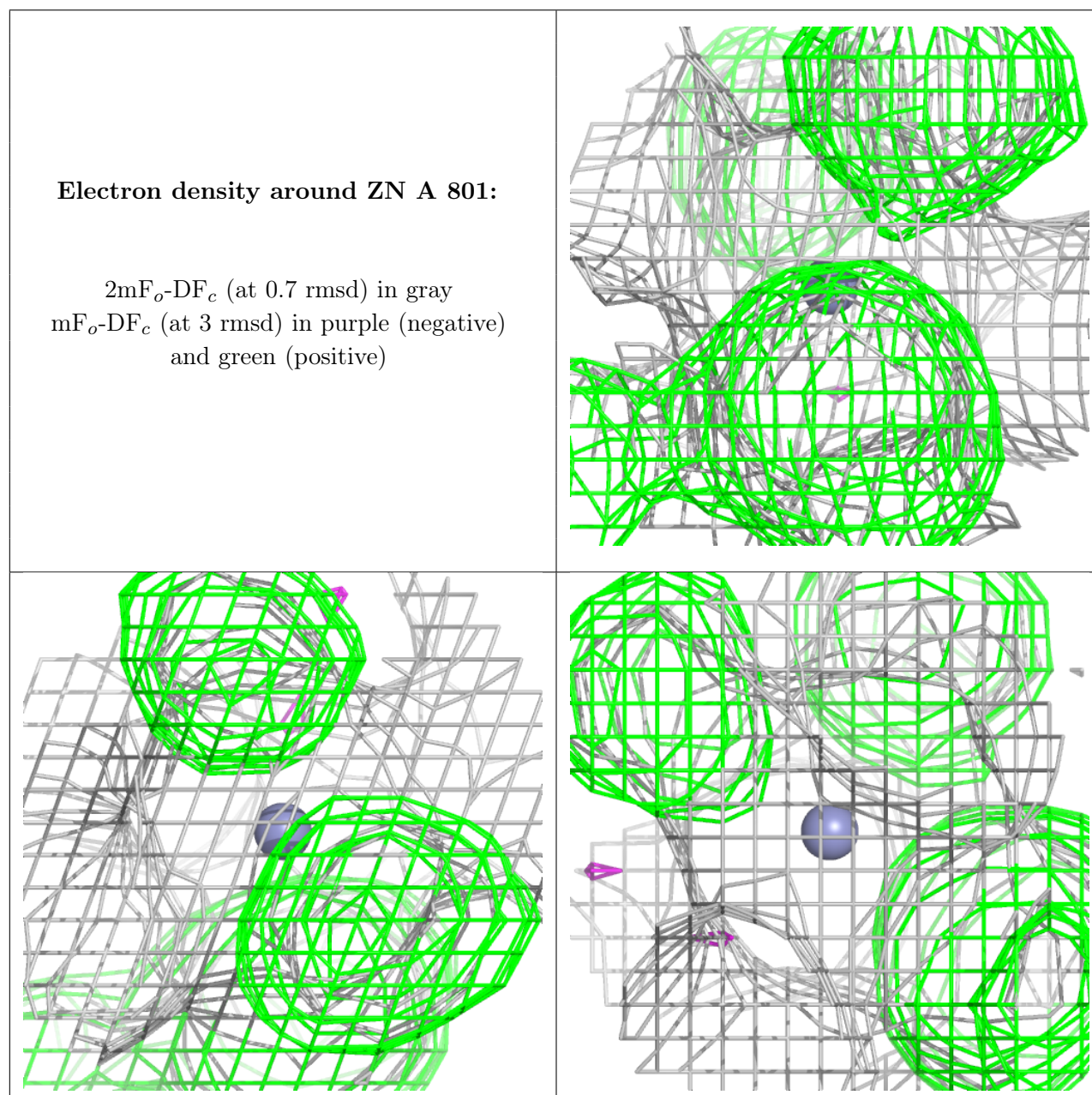
There are no oligosaccharides in this entry.

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.

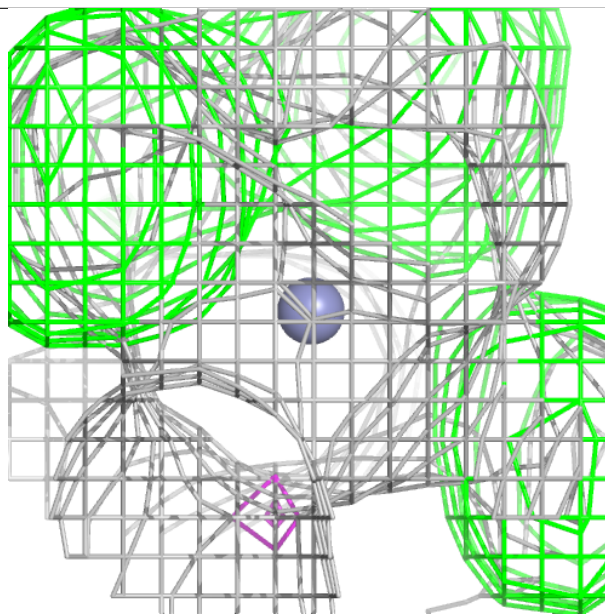
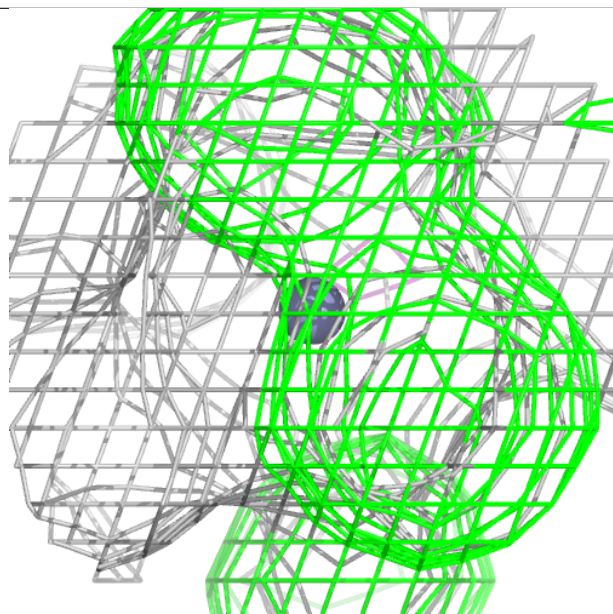
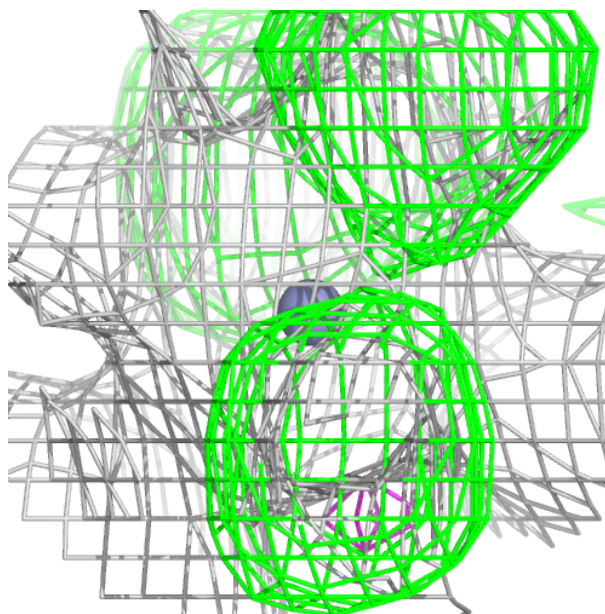
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	ZN	A	801	1/1	0.98	0.04	27,27,27,27	0
2	ZN	B	801	1/1	0.98	0.04	28,28,28,28	0

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 ZN B 801:

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



6.5 Other polymers ⓘ

There are no such residues in this entry.