



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 18, 2026 – 07:23 AM UTC

PDB ID : 7DH7 / pdb_00007dh7
Title : Crystal structure of apo XcZur
Authors : Liu, F.M.; Su, Z.H.; Chen, P.; Tian, X.L.; Wu, L.J.; Tang, D.J.; Li, P.F.;
Deng, H.T.; Tang, J.L.; Ming, Z.H.
Deposited on : 2020-11-13
Resolution : 2.20 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

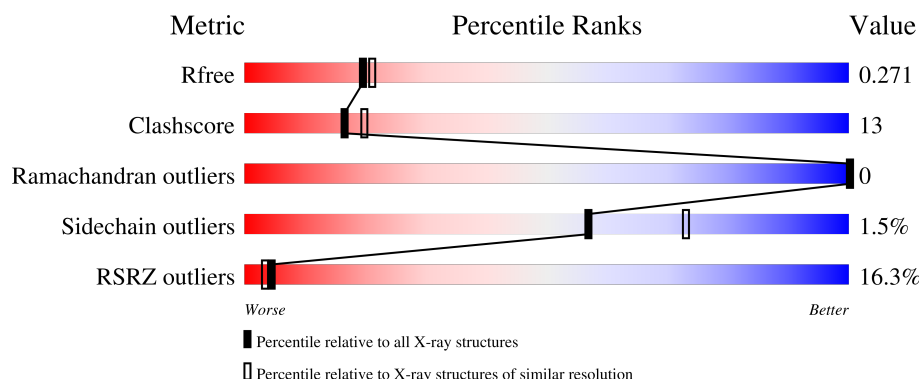
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.20 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	6164 (2.20-2.20)
Clashscore	190562	6851 (2.20-2.20)
Ramachandran outliers	187476	6768 (2.20-2.20)
Sidechain outliers	187428	6769 (2.20-2.20)
RSRZ outliers	180081	6166 (2.20-2.20)

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	170	15% 71%16%13%
1	B	170	22% 72%15%13%
1	C	170	9% 71%18%12%
1	D	170	11% 67%19%..12%

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 4914 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 Transcriptional regulator fur family.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	148	Total	C	N	O	S	0	0	0
			1129	706	210	206	7			
1	B	148	Total	C	N	O	S	0	0	0
			1131	708	210	206	7			
1	C	150	Total	C	N	O	S	0	0	0
			1146	717	214	208	7			
1	D	150	Total	C	N	O	S	0	0	0
			1140	713	212	208	7			

There are 56 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	2	MET	-	initiating methionine	UNP Q4UWS5
A	3	GLY	-	expression tag	UNP Q4UWS5
A	4	SER	-	expression tag	UNP Q4UWS5
A	5	SER	-	expression tag	UNP Q4UWS5
A	6	HIS	-	expression tag	UNP Q4UWS5
A	7	HIS	-	expression tag	UNP Q4UWS5
A	8	HIS	-	expression tag	UNP Q4UWS5
A	9	HIS	-	expression tag	UNP Q4UWS5
A	10	HIS	-	expression tag	UNP Q4UWS5
A	11	HIS	-	expression tag	UNP Q4UWS5
A	12	SER	-	expression tag	UNP Q4UWS5
A	13	GLN	-	expression tag	UNP Q4UWS5
A	14	GLY	-	expression tag	UNP Q4UWS5
A	15	SER	-	expression tag	UNP Q4UWS5
B	2	MET	-	initiating methionine	UNP Q4UWS5
B	3	GLY	-	expression tag	UNP Q4UWS5
B	4	SER	-	expression tag	UNP Q4UWS5
B	5	SER	-	expression tag	UNP Q4UWS5
B	6	HIS	-	expression tag	UNP Q4UWS5
B	7	HIS	-	expression tag	UNP Q4UWS5
B	8	HIS	-	expression tag	UNP Q4UWS5

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Chain	Residue	Modelled	Actual	Comment	Reference
B	9	HIS	-	expression tag	UNP Q4UWS5
B	10	HIS	-	expression tag	UNP Q4UWS5
B	11	HIS	-	expression tag	UNP Q4UWS5
B	12	SER	-	expression tag	UNP Q4UWS5
B	13	GLN	-	expression tag	UNP Q4UWS5
B	14	GLY	-	expression tag	UNP Q4UWS5
B	15	SER	-	expression tag	UNP Q4UWS5
C	2	MET	-	initiating methionine	UNP Q4UWS5
C	3	GLY	-	expression tag	UNP Q4UWS5
C	4	SER	-	expression tag	UNP Q4UWS5
C	5	SER	-	expression tag	UNP Q4UWS5
C	6	HIS	-	expression tag	UNP Q4UWS5
C	7	HIS	-	expression tag	UNP Q4UWS5
C	8	HIS	-	expression tag	UNP Q4UWS5
C	9	HIS	-	expression tag	UNP Q4UWS5
C	10	HIS	-	expression tag	UNP Q4UWS5
C	11	HIS	-	expression tag	UNP Q4UWS5
C	12	SER	-	expression tag	UNP Q4UWS5
C	13	GLN	-	expression tag	UNP Q4UWS5
C	14	GLY	-	expression tag	UNP Q4UWS5
C	15	SER	-	expression tag	UNP Q4UWS5
D	2	MET	-	initiating methionine	UNP Q4UWS5
D	3	GLY	-	expression tag	UNP Q4UWS5
D	4	SER	-	expression tag	UNP Q4UWS5
D	5	SER	-	expression tag	UNP Q4UWS5
D	6	HIS	-	expression tag	UNP Q4UWS5
D	7	HIS	-	expression tag	UNP Q4UWS5
D	8	HIS	-	expression tag	UNP Q4UWS5
D	9	HIS	-	expression tag	UNP Q4UWS5
D	10	HIS	-	expression tag	UNP Q4UWS5
D	11	HIS	-	expression tag	UNP Q4UWS5
D	12	SER	-	expression tag	UNP Q4UWS5
D	13	GLN	-	expression tag	UNP Q4UWS5
D	14	GLY	-	expression tag	UNP Q4UWS5
D	15	SER	-	expression tag	UNP Q4UWS5

- 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 1 1	0	0

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

- Molecule 3 is PHOSPHATE ION (CCD ID: PO4) (formula: O₄P).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	O	P	0	0
			5	4	1		
3	B	1	Total	O	P	0	0
			5	4	1		
3	C	1	Total	O	P	0	0
			5	4	1		
3	D	1	Total	O	P	0	0
			5	4	1		
3	D	1	Total	O	P	0	0
			5	4	1		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	99	Total	O	0	0
			99	99		

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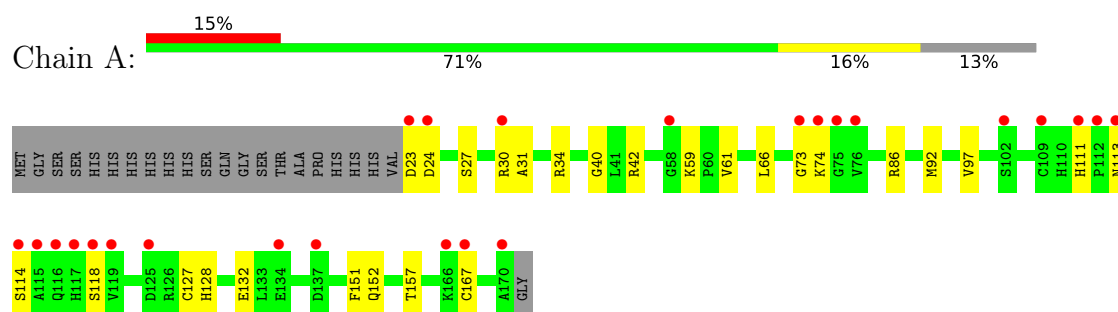
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	B	88	Total 88	O 88	0	0
4	C	81	Total 81	O 81	0	0
4	D	71	Total 71	O 71	0	0

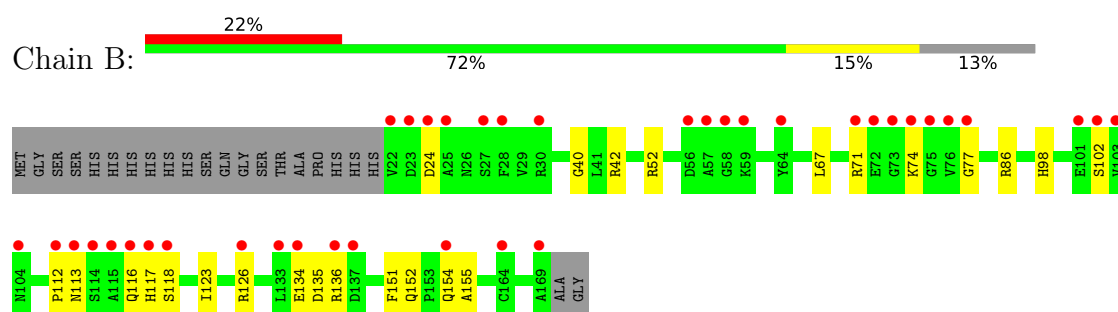
3 Residue-property plots

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.

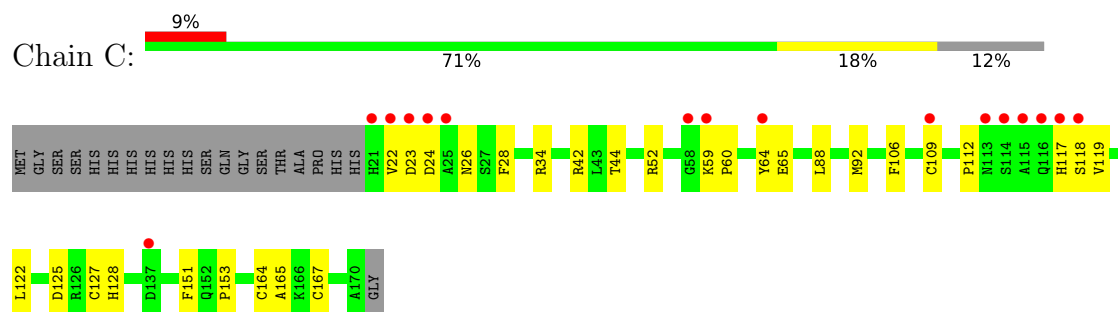
- Molecule 1: Transcriptional regulator fur family



- Molecule 1: Transcriptional regulator fur family

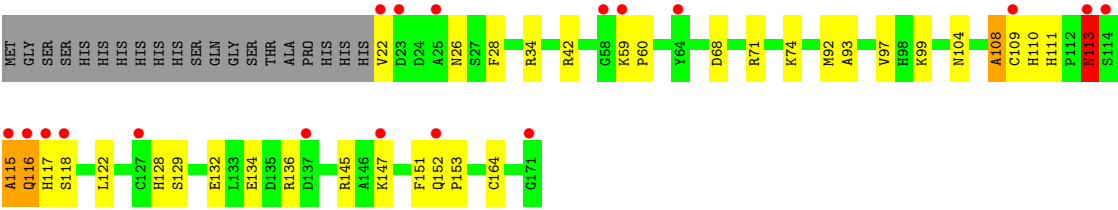


- Molecule 1: Transcriptional regulator fur family



- Molecule 1: Transcriptional regulator fur family





4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	153.39Å 77.87Å 73.06Å 90.00° 103.07° 90.00°	Depositor
Resolution (Å)	38.94 – 2.20 38.94 – 2.20	Depositor EDS
% Data completeness (in resolution range)	99.5 (38.94-2.20) 99.5 (38.94-2.20)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.58 (at 2.20Å)	Xtriage
Refinement program	PHENIX 1.18rc1_3777	Depositor
R, R_{free}	0.219 , 0.276 0.219 , 0.271	Depositor DCC
R_{free} test set	2029 reflections (4.76%)	wwPDB-VP
Wilson B-factor (Å ²)	22.0	Xtriage
Anisotropy	0.000	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 60.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	4914	wwPDB-VP
Average B, all atoms (Å ²)	40.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 16.14% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: PO4, 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	0.46	0/1151	0.67	0/1561
1	B	0.54	0/1153	0.77	2/1564 (0.1%)
1	C	0.66	0/1169	0.87	2/1586 (0.1%)
1	D	0.62	0/1162	0.84	5/1576 (0.3%)
All	All	0.58	0/4635	0.79	9/6287 (0.1%)

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	B	0	1

There are no bond length outliers.

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	125	ASP	CA-CB-CG	5.67	118.27	112.60
1	D	113	ASN	CB-CA-C	5.49	120.59	112.03
1	C	112	PRO	N-CA-C	5.45	119.65	111.14
1	D	115	ALA	CB-CA-C	-5.28	110.47	116.54
1	D	116	GLN	N-CA-C	-5.25	107.42	113.88
1	B	71	ARG	CA-C-N	-5.07	115.64	123.14
1	B	71	ARG	C-N-CA	-5.07	115.64	123.14
1	D	108	ALA	CA-C-N	5.01	128.19	120.82
1	D	108	ALA	C-N-CA	5.01	128.19	120.82

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	113	ASN	Peptide

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	1129	0	1116	36	0
1	B	1131	0	1120	34	0
1	C	1146	0	1132	25	0
1	D	1140	0	1129	43	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
2	C	1	0	0	0	0
2	D	1	0	0	0	0
3	A	5	0	0	0	0
3	B	5	0	0	0	0
3	C	5	0	0	0	0
3	D	10	0	0	0	0
4	A	99	0	0	8	0
4	B	88	0	0	9	0
4	C	81	0	0	1	0
4	D	71	0	0	3	0
All	All	4914	0	4497	117	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 13.

All (117) 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:74:LYS:HA	1:A:74:LYS:HE2	1.36	1.06
1:B:136:ARG:NH1	4:B:302:HOH:O	2.02	0.93
1:D:116:GLN:OE1	4:D:301:HOH:O	1.91	0.88
1:D:147:LYS:HE3	1:D:152:GLN:HE22	1.39	0.88
1:B:151:PHE:CD1	1:D:122:LEU:HD21	2.09	0.86
1:B:77:GLY:HA2	1:C:34:ARG:HH21	1.42	0.84
1:B:151:PHE:CD1	1:D:122:LEU:CD2	2.66	0.79

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:154:GLN:NE2	4:B:303:HOH:O	2.13	0.77
1:A:27:SER:OG	4:A:301:HOH:O	2.03	0.76
1:C:59:LYS:HG3	1:C:60:PRO:HD2	1.69	0.74
1:B:151:PHE:CE1	1:D:122:LEU:HD23	2.23	0.73
1:B:116:GLN:O	1:B:117:HIS:ND1	2.22	0.72
1:A:23:ASP:N	4:A:302:HOH:O	2.22	0.72
1:D:92:MET:HE2	1:D:128:HIS:HB2	1.72	0.72
1:B:42:ARG:HG3	1:C:42:ARG:HG3	1.71	0.71
1:D:117:HIS:CD2	1:D:132:GLU:OE2	2.44	0.71
1:A:74:LYS:HE2	1:A:74:LYS:CA	2.19	0.71
1:B:151:PHE:CE1	1:D:122:LEU:CD2	2.74	0.70
1:A:73:GLY:O	1:A:74:LYS:HD2	1.93	0.68
1:D:68:ASP:OD1	4:D:302:HOH:O	2.13	0.66
1:D:110:HIS:ND1	1:D:115:ALA:HB1	2.10	0.66
1:D:147:LYS:HE3	1:D:152:GLN:NE2	2.10	0.66
1:A:74:LYS:HA	1:A:74:LYS:CE	2.07	0.65
1:D:110:HIS:HA	1:D:115:ALA:HB1	1.79	0.65
1:A:24:ASP:O	1:A:27:SER:OG	2.15	0.64
1:B:74:LYS:HE3	1:C:22:VAL:HG21	1.79	0.64
1:D:59:LYS:HD3	1:D:60:PRO:O	1.97	0.64
1:B:77:GLY:HA2	1:C:34:ARG:NH2	2.13	0.63
1:D:108:ALA:O	1:D:116:GLN:HG3	1.99	0.63
1:D:116:GLN:NE2	4:D:304:HOH:O	2.29	0.63
1:B:116:GLN:H	1:B:116:GLN:CD	2.08	0.62
1:D:136:ARG:HD3	1:D:136:ARG:H	1.64	0.62
1:A:42:ARG:HG3	1:D:42:ARG:HG3	1.82	0.61
1:B:74:LYS:HE3	1:C:22:VAL:CG2	2.31	0.61
1:A:74:LYS:HD2	4:A:321:HOH:O	2.01	0.61
1:A:42:ARG:NH1	1:B:40:GLY:O	2.31	0.60
1:B:117:HIS:O	1:B:117:HIS:CG	2.54	0.60
1:D:108:ALA:HB3	1:D:116:GLN:HG2	1.85	0.59
1:A:40:GLY:HA3	4:A:341:HOH:O	2.03	0.59
1:D:22:VAL:HG12	1:D:26:ASN:ND2	2.19	0.57
1:B:136:ARG:NH2	4:B:309:HOH:O	2.38	0.57
1:D:108:ALA:N	1:D:116:GLN:HG3	2.19	0.57
1:A:127:CYS:HB3	1:A:167:CYS:SG	2.46	0.56
1:A:30:ARG:HG3	1:A:31:ALA:N	2.20	0.56
1:A:30:ARG:NH2	1:A:34:ARG:HH11	2.04	0.55
1:C:23:ASP:CG	1:C:24:ASP:H	2.15	0.55
1:D:111:HIS:H	1:D:115:ALA:HA	1.71	0.55
1:B:136:ARG:NH2	4:B:304:HOH:O	2.39	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:135:ASP:HB2	1:D:145:ARG:HD3	1.89	0.54
1:B:118:SER:N	4:B:312:HOH:O	2.41	0.54
1:B:102:SER:N	4:B:313:HOH:O	2.41	0.54
1:B:151:PHE:CD1	1:D:122:LEU:HD23	2.42	0.54
1:B:52:ARG:HD2	4:B:372:HOH:O	2.08	0.54
1:A:132:GLU:OE1	1:A:132:GLU:N	2.42	0.53
1:D:28:PHE:HE1	1:D:109:CYS:SG	2.33	0.52
1:C:22:VAL:HB	1:C:26:ASN:HB3	1.90	0.52
1:D:92:MET:HA	1:D:97:VAL:O	2.11	0.51
1:A:132:GLU:H	1:A:132:GLU:CD	2.18	0.51
1:B:152:GLN:NE2	4:B:301:HOH:O	1.96	0.51
1:A:152:GLN:HG3	1:C:165:ALA:HA	1.93	0.50
1:D:113:ASN:HD22	1:D:113:ASN:N	2.10	0.50
1:C:28:PHE:HE1	1:C:109:CYS:HG	1.57	0.50
1:A:27:SER:HA	1:A:30:ARG:HG2	1.95	0.49
1:A:61:VAL:HG11	1:A:66:LEU:HG	1.92	0.49
1:C:127:CYS:HB3	1:C:167:CYS:SG	2.52	0.49
1:C:28:PHE:HE1	1:C:109:CYS:SG	2.35	0.49
1:D:113:ASN:N	1:D:113:ASN:ND2	2.60	0.48
1:A:152:GLN:NE2	4:A:312:HOH:O	2.44	0.48
1:A:118:SER:HB3	1:A:157:THR:OG1	2.13	0.48
1:A:74:LYS:CD	4:A:321:HOH:O	2.60	0.48
1:C:92:MET:HE2	1:C:128:HIS:HB2	1.96	0.47
1:D:136:ARG:H	1:D:136:ARG:CD	2.26	0.47
1:A:86:ARG:HG2	1:B:86:ARG:NE	2.29	0.47
1:C:26:ASN:OD1	1:C:52:ARG:NH2	2.48	0.47
1:A:111:HIS:NE2	4:A:301:HOH:O	2.36	0.46
1:B:126:ARG:HH21	1:D:34:ARG:NH2	2.13	0.46
1:B:67:LEU:HD12	1:B:67:LEU:O	2.16	0.46
1:D:151:PHE:O	1:D:153:PRO:HD3	2.16	0.45
1:D:136:ARG:HD3	1:D:136:ARG:N	2.30	0.45
1:D:108:ALA:O	1:D:116:GLN:CG	2.64	0.45
1:C:23:ASP:OD1	1:C:24:ASP:N	2.44	0.44
1:A:30:ARG:HH22	1:A:34:ARG:HH11	1.64	0.44
1:B:117:HIS:O	1:B:117:HIS:CD2	2.70	0.44
1:A:111:HIS:O	1:A:113:ASN:N	2.50	0.44
1:B:134:GLU:HB3	4:B:304:HOH:O	2.17	0.44
1:A:86:ARG:NE	1:B:86:ARG:HG2	2.33	0.44
1:A:24:ASP:O	4:A:301:HOH:O	2.21	0.43
1:B:98:HIS:CE1	1:B:123:ILE:HD13	2.52	0.43
1:D:93:ALA:HA	1:D:128:HIS:CE1	2.53	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:24:ASP:N	1:B:24:ASP:OD1	2.50	0.43
1:D:104:ASN:O	1:D:104:ASN:ND2	2.50	0.43
1:B:126:ARG:NH2	1:D:34:ARG:NH2	2.66	0.43
1:C:26:ASN:OD1	1:C:52:ARG:NH1	2.52	0.43
1:C:151:PHE:O	1:C:153:PRO:HD3	2.19	0.42
1:A:113:ASN:CG	1:A:114:SER:H	2.23	0.42
1:D:108:ALA:H	1:D:116:GLN:NE2	2.17	0.42
1:A:59:LYS:HE3	1:A:59:LYS:HB2	1.87	0.42
1:D:71:ARG:O	1:D:74:LYS:O	2.38	0.42
1:D:147:LYS:HA	1:D:151:PHE:O	2.19	0.42
1:A:74:LYS:CA	1:A:74:LYS:CE	2.83	0.42
1:C:42:ARG:HH21	1:C:44:THR:HG22	1.85	0.42
1:C:28:PHE:CE1	1:C:109:CYS:SG	3.08	0.42
1:C:64:TYR:CE1	1:C:65:GLU:HG3	2.55	0.42
1:D:108:ALA:N	1:D:116:GLN:CG	2.83	0.42
1:A:92:MET:HA	1:A:97:VAL:O	2.20	0.42
1:A:151:PHE:CE1	1:C:122:LEU:HG	2.54	0.41
1:C:88:LEU:HD21	1:C:106:PHE:CZ	2.54	0.41
1:A:92:MET:HE2	1:A:128:HIS:HB2	2.02	0.41
1:D:134:GLU:O	1:D:136:ARG:HD3	2.19	0.41
1:A:40:GLY:O	1:B:42:ARG:NH1	2.45	0.41
1:D:28:PHE:CE1	1:D:109:CYS:SG	3.09	0.41
1:B:112:PRO:HA	1:B:155:ALA:O	2.20	0.41
1:D:108:ALA:H	1:D:116:GLN:HE21	1.67	0.41
1:A:152:GLN:CG	1:C:165:ALA:HA	2.51	0.41
1:C:34:ARG:NH1	4:C:316:HOH:O	2.54	0.40
1:D:99:LYS:NZ	1:D:104:ASN:OD1	2.52	0.40
1:C:117:HIS:C	1:C:117:HIS:HD1	2.30	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	146/170 (86%)	139 (95%)	7 (5%)	0	100	100
1	B	146/170 (86%)	140 (96%)	6 (4%)	0	100	100
1	C	148/170 (87%)	144 (97%)	4 (3%)	0	100	100
1	D	148/170 (87%)	143 (97%)	5 (3%)	0	100	100
All	All	588/680 (86%)	566 (96%)	22 (4%)	0	100	100

There are no Ramachandran outliers to report.

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	117/135 (87%)	117 (100%)	0	100	100
1	B	118/135 (87%)	118 (100%)	0	100	100
1	C	119/135 (88%)	116 (98%)	3 (2%)	42	56
1	D	118/135 (87%)	114 (97%)	4 (3%)	32	44
All	All	472/540 (87%)	465 (98%)	7 (2%)	57	73

All (7) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	C	118	SER
1	C	119	VAL
1	C	164	CYS
1	D	113	ASN
1	D	118	SER
1	D	129	SER
1	D	164	CYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	104	ASN

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Mol	Chain	Res	Type
1	B	116	GLN
1	C	152	GLN
1	D	113	ASN
1	D	116	GLN
1	D	117	HIS
1	D	152	GLN
1	D	156	GLN

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 9 ligands modelled in this entry, 4 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	$\# Z > 2$
3	PO4	D	202	-	4,4,4	1.10	0	6,6,6	0.86	0
3	PO4	A	202	-	4,4,4	1.42	0	6,6,6	1.30	0
3	PO4	C	202	-	4,4,4	1.00	0	6,6,6	0.41	0
3	PO4	D	203	-	4,4,4	0.92	0	6,6,6	0.43	0
3	PO4	B	202	-	4,4,4	1.12	0	6,6,6	0.62	0

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	148/170 (87%)	0.90	25 (16%) 4 3	11, 29, 78, 152	0
1	B	148/170 (87%)	1.17	38 (25%) 1 1	12, 36, 78, 130	0
1	C	150/170 (88%)	0.63	16 (10%) 11 8	10, 36, 88, 132	0
1	D	150/170 (88%)	0.74	18 (12%) 9 6	10, 38, 73, 127	0
All	All	596/680 (87%)	0.86	97 (16%) 4 3	10, 36, 82, 152	0

All (97) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	22	VAL	9.6
1	A	113	ASN	8.9
1	B	76	VAL	8.0
1	D	117	HIS	7.8
1	A	117	HIS	7.5
1	B	117	HIS	7.0
1	B	115	ALA	6.5
1	A	73	GLY	6.3
1	A	116	GLN	5.7
1	C	109	CYS	5.7
1	A	102	SER	5.5
1	A	115	ALA	5.4
1	B	73	GLY	5.2
1	C	23	ASP	5.1
1	B	74	LYS	5.1
1	D	58	GLY	5.0
1	C	113	ASN	5.0
1	B	114	SER	4.9
1	A	166	LYS	4.6
1	A	170	ALA	4.5
1	C	117	HIS	4.5

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Mol	Chain	Res	Type	RSRZ
1	B	22	VAL	4.4
1	B	113	ASN	4.4
1	B	58	GLY	4.4
1	D	114	SER	4.4
1	A	74	LYS	4.4
1	D	109	CYS	4.3
1	A	75	GLY	4.3
1	D	116	GLN	4.2
1	A	119	VAL	4.2
1	B	72	GLU	4.2
1	A	118	SER	4.1
1	D	22	VAL	4.1
1	A	23	ASP	4.1
1	A	24	ASP	4.1
1	C	115	ALA	4.0
1	D	115	ALA	4.0
1	D	23	ASP	3.9
1	A	114	SER	3.9
1	B	118	SER	3.8
1	C	21	HIS	3.6
1	B	59	LYS	3.6
1	B	154	GLN	3.6
1	B	23	ASP	3.6
1	D	127	CYS	3.6
1	A	109	CYS	3.5
1	C	59	LYS	3.5
1	B	77	GLY	3.5
1	C	114	SER	3.5
1	A	137	ASP	3.4
1	B	134	GLU	3.3
1	B	27	SER	3.2
1	D	64	TYR	3.2
1	B	102	SER	3.2
1	B	30	ARG	3.2
1	A	76	VAL	3.1
1	A	167	CYS	2.9
1	C	58	GLY	2.9
1	B	133	LEU	2.9
1	D	113	ASN	2.9
1	B	116	GLN	2.8
1	D	152	GLN	2.7
1	B	103	VAL	2.7

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Mol	Chain	Res	Type	RSRZ
1	D	137	ASP	2.7
1	A	134	GLU	2.7
1	D	59	LYS	2.7
1	D	25	ALA	2.7
1	C	116	GLN	2.7
1	D	171	GLY	2.7
1	A	30	ARG	2.7
1	B	24	ASP	2.6
1	C	137	ASP	2.6
1	B	112	PRO	2.6
1	C	118	SER	2.6
1	B	71	ARG	2.5
1	B	169	ALA	2.5
1	B	75	GLY	2.5
1	B	57	ALA	2.4
1	B	136	ARG	2.3
1	A	112	PRO	2.3
1	B	101	GLU	2.3
1	B	137	ASP	2.3
1	A	125	ASP	2.3
1	B	104	ASN	2.3
1	C	24	ASP	2.3
1	A	58	GLY	2.3
1	B	126	ARG	2.2
1	C	25	ALA	2.2
1	D	118	SER	2.2
1	B	64	TYR	2.1
1	C	64	TYR	2.1
1	D	147	LYS	2.1
1	B	164	CYS	2.1
1	B	56	ASP	2.1
1	B	25	ALA	2.1
1	B	28	PHE	2.1
1	A	111	HIS	2.0

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

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](#)

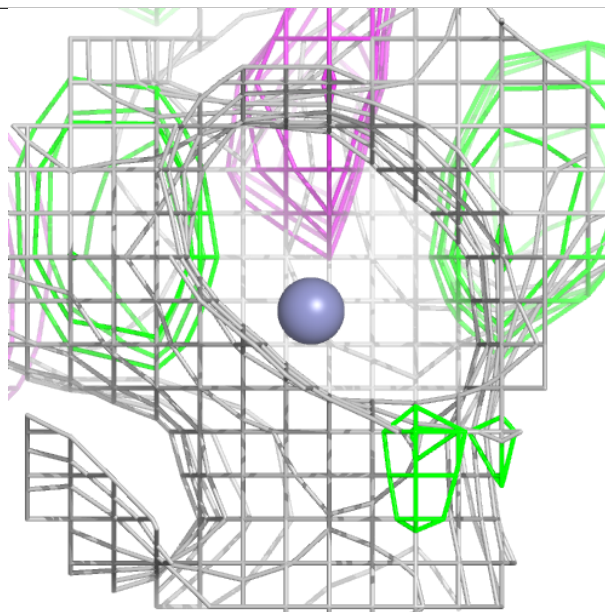
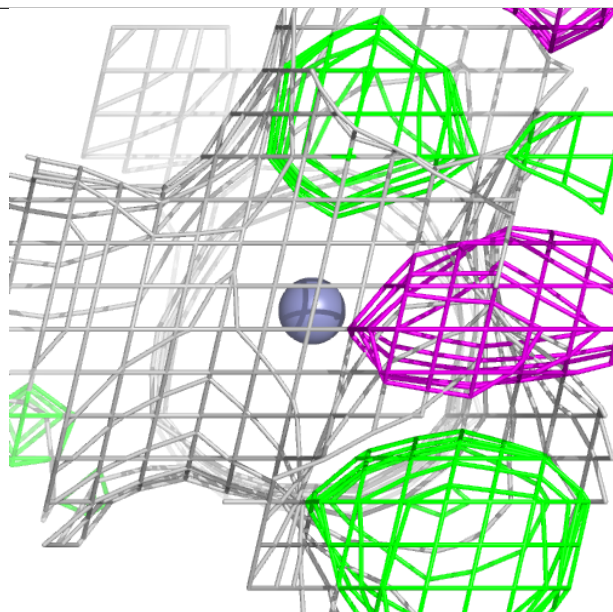
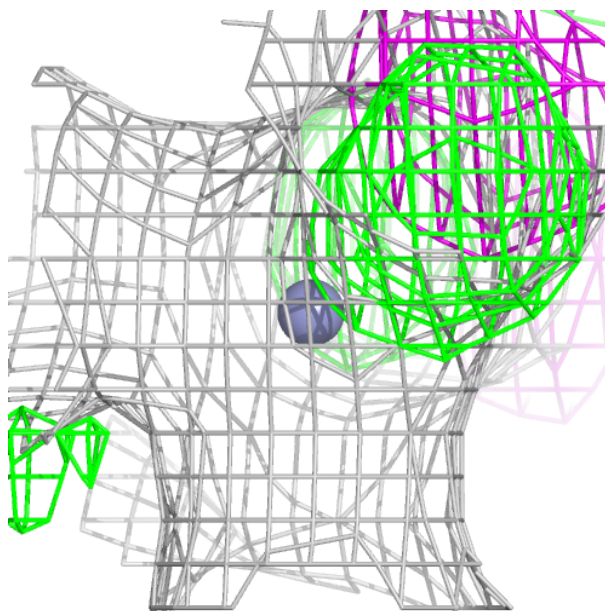
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
3	PO4	C	202	5/5	0.87	0.13	74,78,84,84	0
3	PO4	D	203	5/5	0.90	0.11	62,67,71,75	0
3	PO4	D	202	5/5	0.92	0.13	38,42,46,51	0
3	PO4	B	202	5/5	0.93	0.13	39,40,43,54	0
3	PO4	A	202	5/5	0.99	0.06	15,25,26,27	0
2	ZN	A	201	1/1	0.99	0.02	29,29,29,29	0
2	ZN	B	201	1/1	0.99	0.03	39,39,39,39	0
2	ZN	C	201	1/1	0.99	0.04	28,28,28,28	0
2	ZN	D	201	1/1	0.99	0.07	34,34,34,34	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.

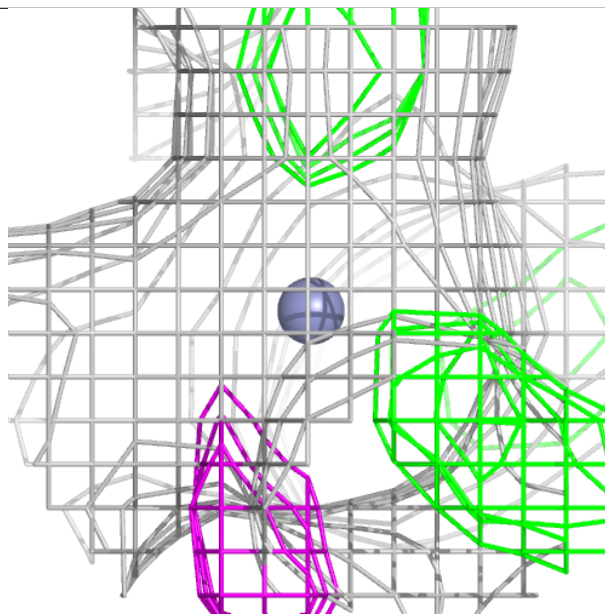
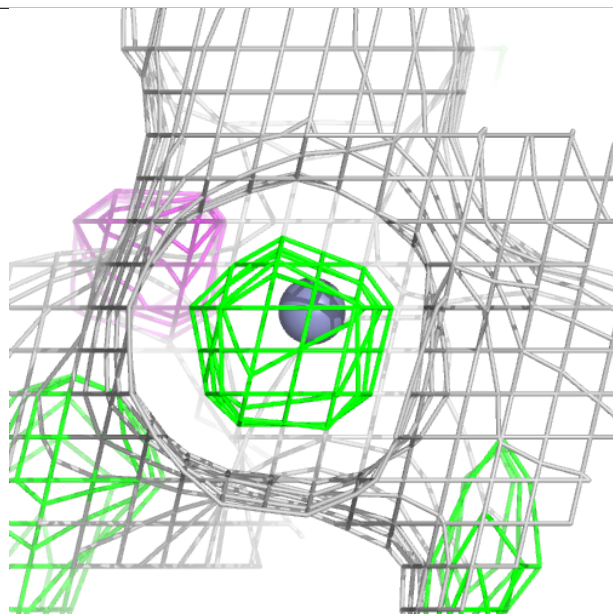
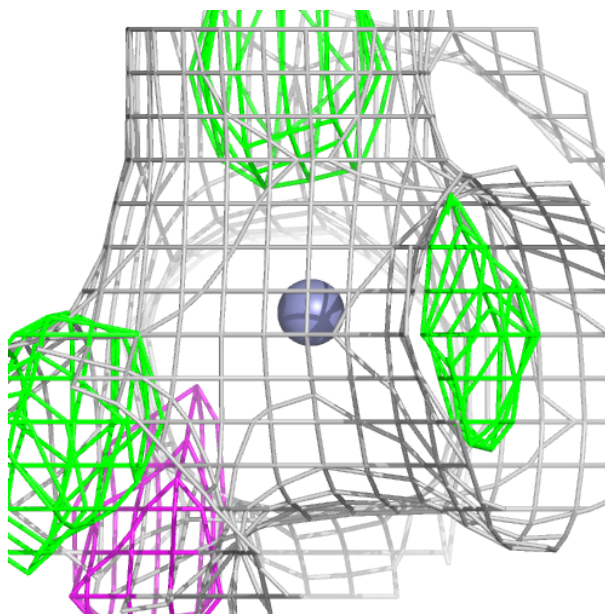
Electron density around ZN A 201:

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



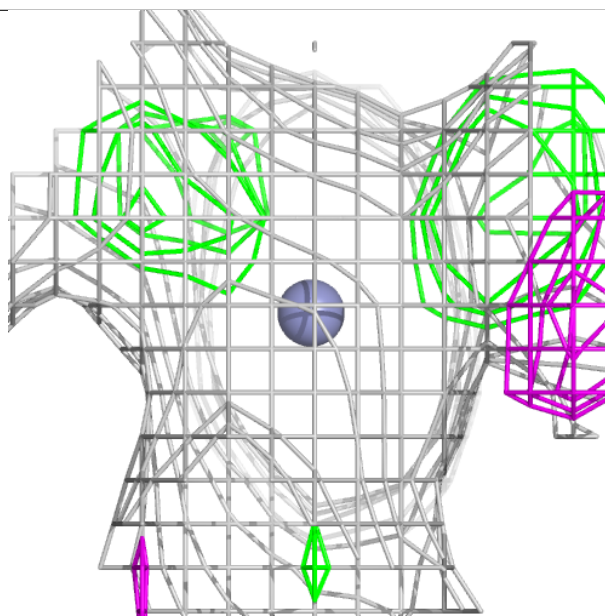
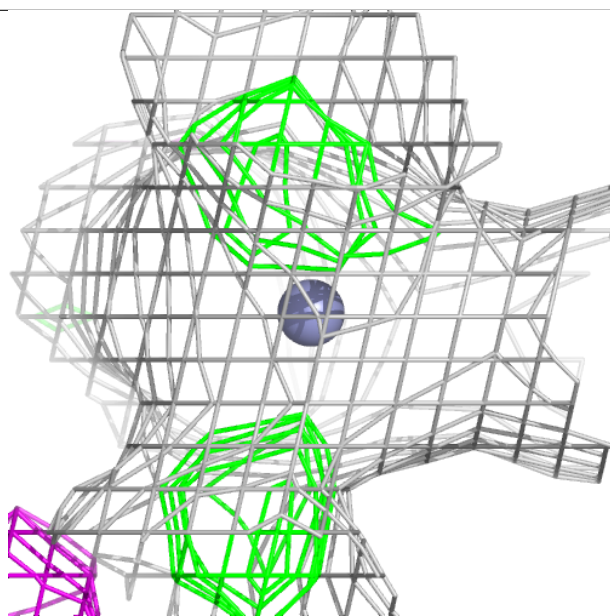
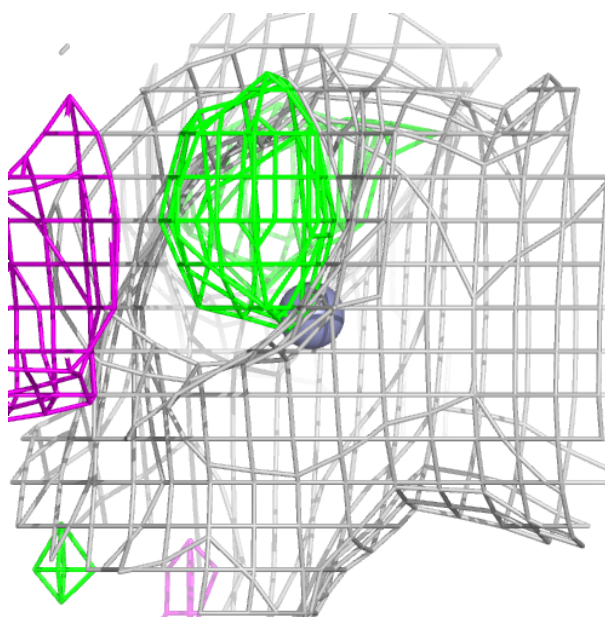
Electron density around ZN B 201:

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



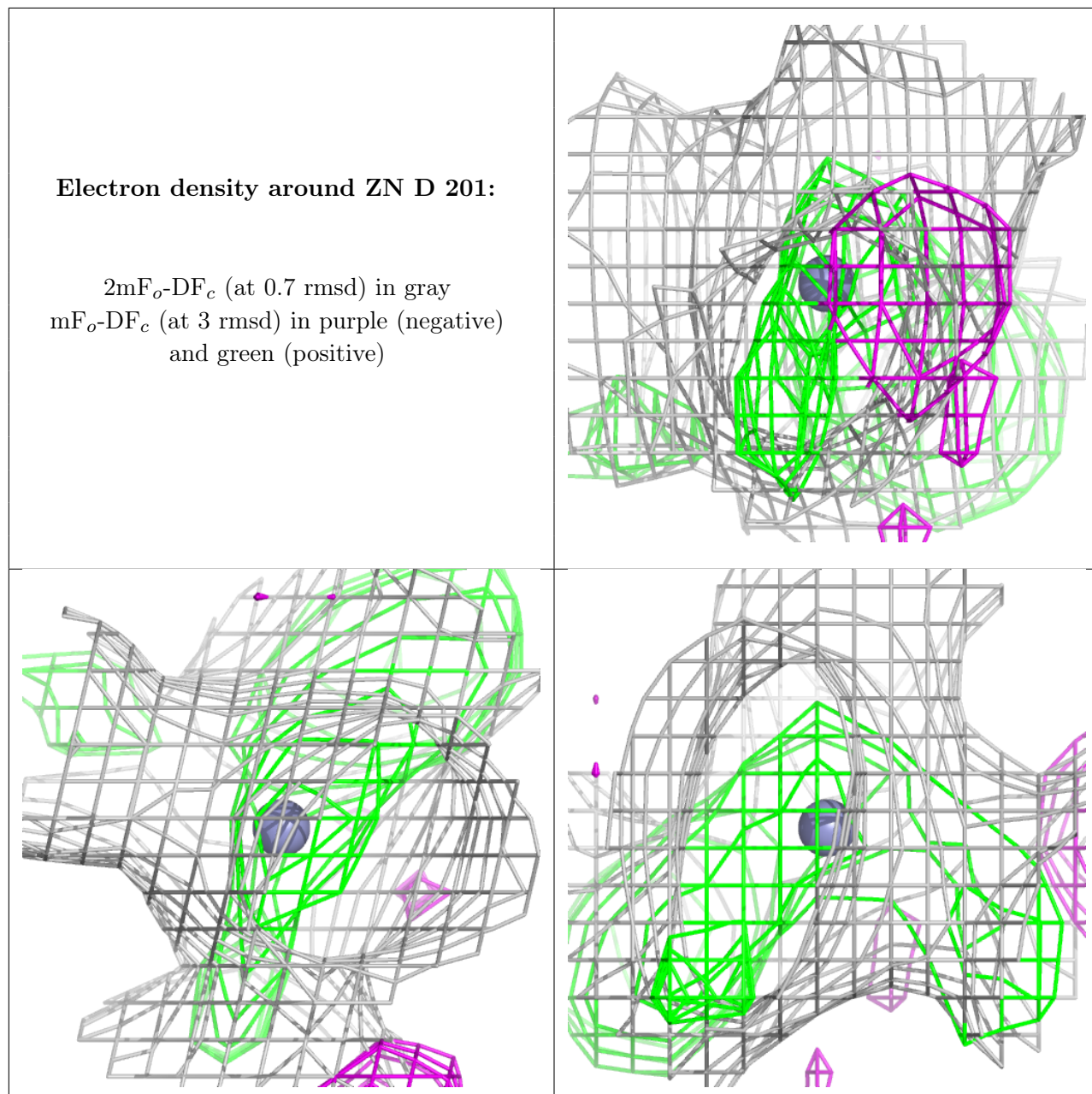
Electron density around ZN C 201:

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



Electron density around ZN D 201:

$2mF_o-DF_c$ (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.