



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 10:49 PM UTC

PDB ID : 7WI1 / pdb_00007wi1
Title : The mutant variant of PNGM-1, H93 was substituted for alanine to study metal coordination
Authors : Park, Y.S.; Kang, L.W.; Lee, J.H.
Deposited on : 2022-01-01
Resolution : 1.61 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

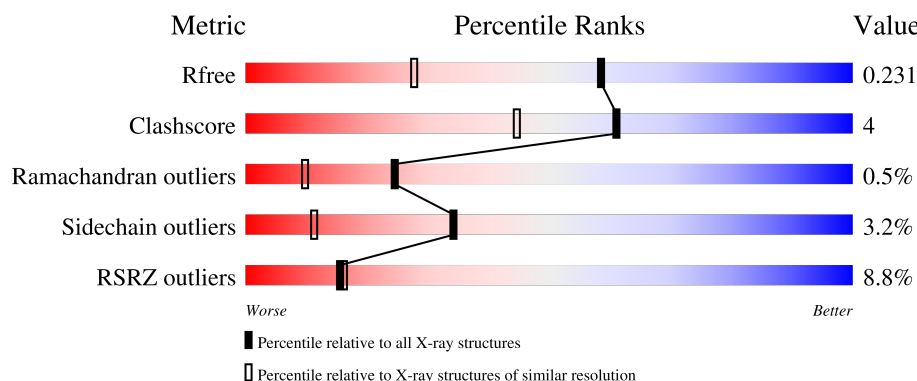
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.61 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	6728 (1.64-1.60)
Clashscore	190562	7023 (1.64-1.60)
Ramachandran outliers	187476	6898 (1.64-1.60)
Sidechain outliers	187428	6896 (1.64-1.60)
RSRZ outliers	180081	6727 (1.64-1.60)

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	372	7% 88% 10% .
1	B	372	10% 90% 9% .
1	C	372	6% 91% 8% .
1	D	372	7% 94% 5% ..
1	E	372	7% 90% 9% ..

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Mol	Chain	Length	Quality of chain
1	F	372	7% 90% 9% ..
1	G	372	15% 89% 10% .
1	H	372	11% 89% 10% .

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 24542 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 Metallo-beta-lactamase PNGM-1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	372	Total	C	N	O	S	0	0	0
			2928	1850	504	554	20			
1	G	372	Total	C	N	O	S	0	0	0
			2928	1850	504	554	20			
1	C	372	Total	C	N	O	S	0	0	0
			2928	1850	504	554	20			
1	E	372	Total	C	N	O	S	0	0	0
			2928	1850	504	554	20			
1	D	372	Total	C	N	O	S	0	0	0
			2928	1850	504	554	20			
1	F	372	Total	C	N	O	S	0	0	0
			2928	1850	504	554	20			
1	B	372	Total	C	N	O	S	0	0	0
			2928	1850	504	554	20			
1	H	372	Total	C	N	O	S	0	0	0
			2928	1850	504	554	20			

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	93	ALA	HIS	engineered mutation	UNP A0A2U8UYM6
G	93	ALA	HIS	engineered mutation	UNP A0A2U8UYM6
C	93	ALA	HIS	engineered mutation	UNP A0A2U8UYM6
E	93	ALA	HIS	engineered mutation	UNP A0A2U8UYM6
D	93	ALA	HIS	engineered mutation	UNP A0A2U8UYM6
F	93	ALA	HIS	engineered mutation	UNP A0A2U8UYM6
B	93	ALA	HIS	engineered mutation	UNP A0A2U8UYM6
H	93	ALA	HIS	engineered mutation	UNP A0A2U8UYM6

- 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
2	G	1	Total Zn 1 1	0	0
2	C	1	Total Zn 1 1	0	0
2	E	1	Total Zn 1 1	0	0
2	D	1	Total Zn 1 1	0	0
2	F	1	Total Zn 1 1	0	0
2	B	1	Total Zn 1 1	0	0
2	H	1	Total Zn 1 1	0	0

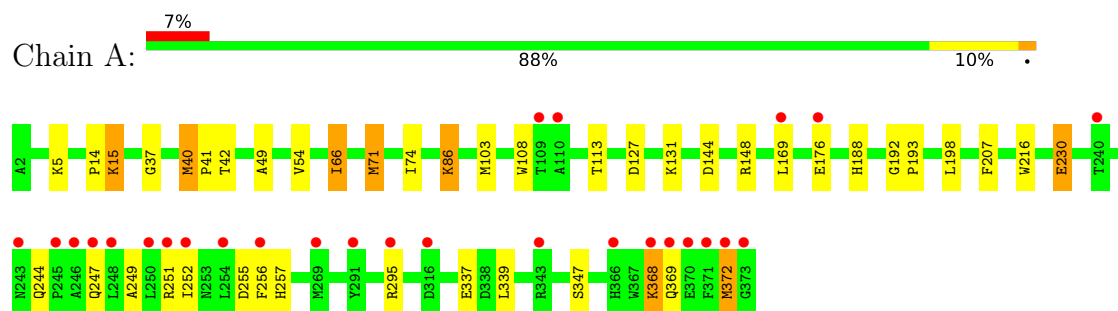
- Molecule 3 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	177	Total O 177 177	0	0
3	G	96	Total O 96 96	0	0
3	C	182	Total O 182 182	0	0
3	E	105	Total O 105 105	0	0
3	D	171	Total O 171 171	0	0
3	F	95	Total O 95 95	0	0
3	B	181	Total O 181 181	0	0
3	H	103	Total O 103 103	0	0

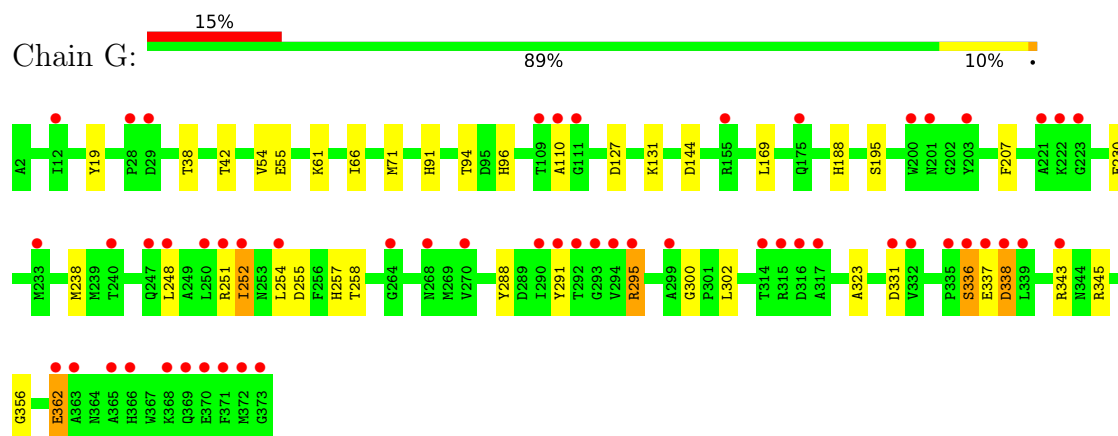
3 Residue-property plots [i](#)

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.

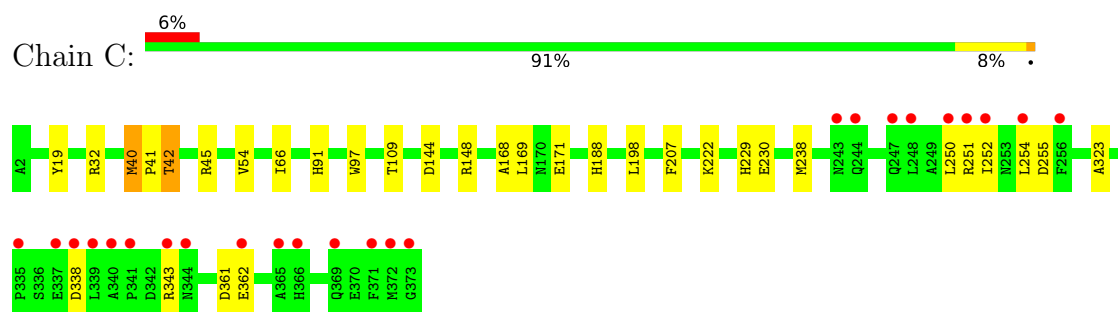
• Molecule 1: Metallo-beta-lactamase PNGM-1



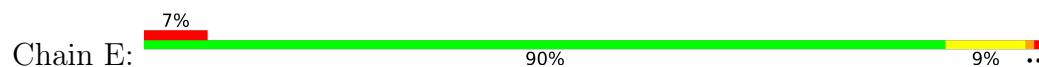
• Molecule 1: Metallo-beta-lactamase PNGM-1

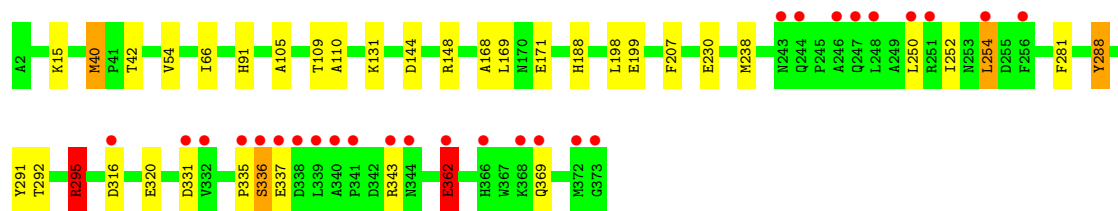


• Molecule 1: Metallo-beta-lactamase PNGM-1

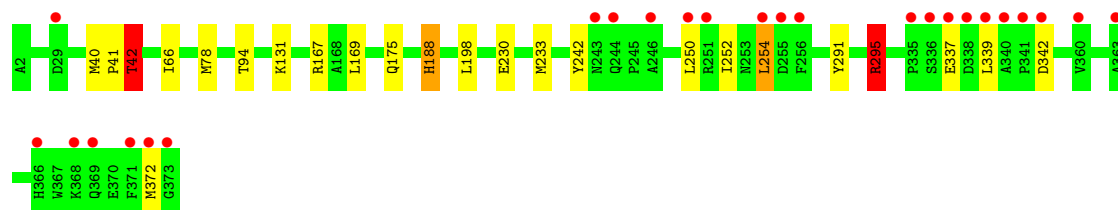


• Molecule 1: Metallo-beta-lactamase PNGM-1

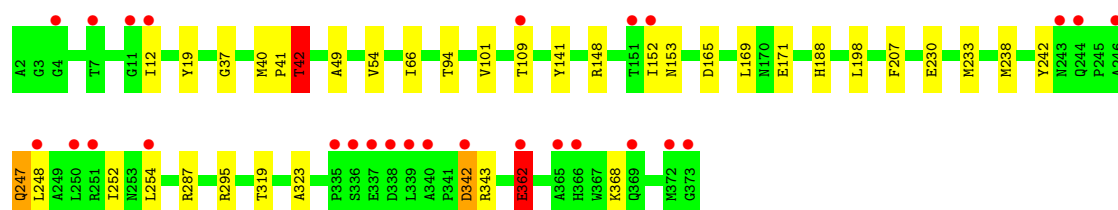
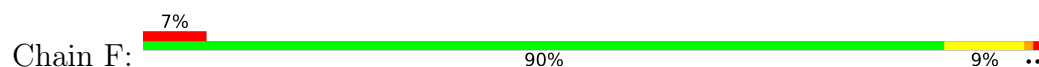




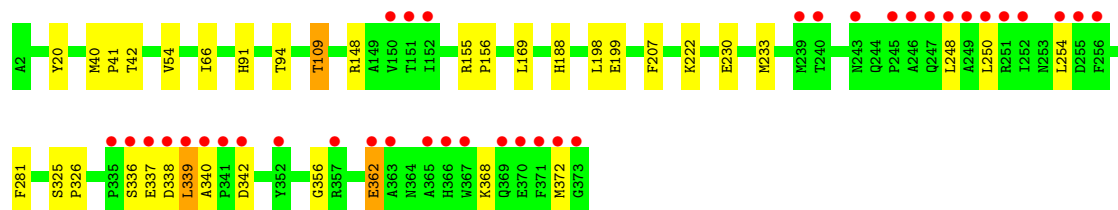
• Molecule 1: Metallo-beta-lactamase PNGM-1



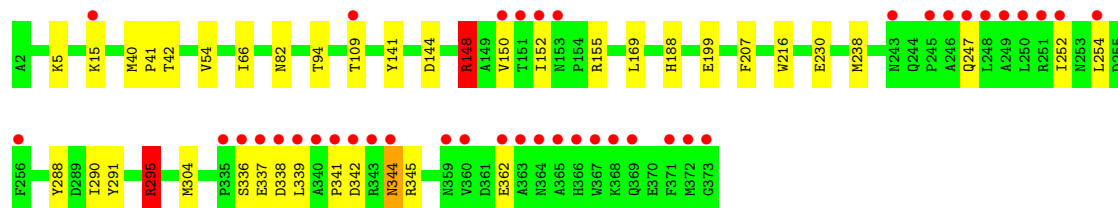
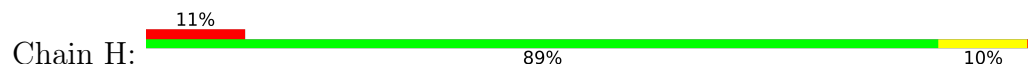
• Molecule 1: Metallo-beta-lactamase PNGM-1



• Molecule 1: Metallo-beta-lactamase PNGM-1



• Molecule 1: Metallo-beta-lactamase PNGM-1



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	121.75Å 82.79Å 164.05Å 90.00° 111.10° 90.00°	Depositor
Resolution (Å)	41.51 – 1.61 41.51 – 1.61	Depositor EDS
% Data completeness (in resolution range)	99.0 (41.51-1.61) 99.0 (41.51-1.61)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.64 (at 1.61Å)	Xtriage
Refinement program	REFMAC 5.8.0267	Depositor
R, R_{free}	0.196 , 0.223 0.205 , 0.231	Depositor DCC
R_{free} test set	19534 reflections (4.98%)	wwPDB-VP
Wilson B-factor (Å ²)	10.3	Xtriage
Anisotropy	0.058	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 32.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.009 for h,-k,-h-l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	24542	wwPDB-VP
Average B, all atoms (Å ²)	14.0	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 82.34 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 2.8594e-07. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

¹Intensities estimated from amplitudes.

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

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: 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	1.16	1/3013 (0.0%)	1.28	2/4107 (0.0%)
1	B	1.11	1/3013 (0.0%)	1.22	2/4107 (0.0%)
1	C	1.08	3/3013 (0.1%)	1.19	3/4107 (0.1%)
1	D	1.12	1/3013 (0.0%)	1.27	5/4107 (0.1%)
1	E	1.10	4/3013 (0.1%)	1.25	10/4107 (0.2%)
1	F	1.10	1/3013 (0.0%)	1.25	5/4107 (0.1%)
1	G	1.15	3/3013 (0.1%)	1.36	8/4107 (0.2%)
1	H	1.11	1/3013 (0.0%)	1.27	8/4107 (0.2%)
All	All	1.12	15/24104 (0.1%)	1.26	43/32856 (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

All (15) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	G	96	HIS	CE1-NE2	8.08	1.40	1.32
1	B	199	GLU	CD-OE1	7.89	1.40	1.25
1	H	345	ARG	CD-NE	-6.81	1.36	1.46
1	E	199	GLU	CD-OE1	6.54	1.37	1.25
1	G	195	SER	C-O	6.39	1.32	1.23
1	F	171	GLU	C-O	6.30	1.31	1.23
1	A	347	SER	C-O	5.73	1.30	1.23
1	E	91	HIS	CE1-NE2	-5.66	1.26	1.32
1	C	45	ARG	C-O	5.46	1.30	1.24
1	E	362	GLU	CD-OE2	5.45	1.35	1.25

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	78	MET	C-O	5.18	1.30	1.23
1	E	320	GLU	C-O	5.12	1.30	1.24
1	G	38	THR	C-O	5.12	1.30	1.24
1	C	229	HIS	CE1-NE2	5.05	1.37	1.32
1	C	97	TRP	N-CA	5.05	1.50	1.46

All (43) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	G	345	ARG	CG-CD-NE	-12.46	84.58	112.00
1	H	345	ARG	CG-CD-NE	-11.03	87.73	112.00
1	F	42	THR	CB-CA-C	9.61	125.63	109.48
1	E	295	ARG	NE-CZ-NH1	9.27	130.77	121.50
1	D	42	THR	CB-CA-C	9.24	125.01	109.48
1	E	295	ARG	NE-CZ-NH2	-8.59	111.47	119.20
1	H	295	ARG	NE-CZ-NH1	8.43	129.93	121.50
1	G	96	HIS	CE1-NE2-CD2	-8.25	100.75	109.00
1	E	295	ARG	CD-NE-CZ	8.21	135.89	124.40
1	H	295	ARG	CD-NE-CZ	7.54	134.96	124.40
1	H	345	ARG	NE-CZ-NH2	-7.52	112.43	119.20
1	B	199	GLU	CB-CG-CD	7.52	125.38	112.60
1	G	71	MET	CG-SD-CE	-7.01	85.48	100.90
1	F	42	THR	N-CA-CB	-6.89	99.15	110.52
1	H	295	ARG	NE-CZ-NH2	-6.56	113.30	119.20
1	E	295	ARG	CB-CG-CD	6.55	126.36	111.30
1	D	295	ARG	CB-CG-CD	6.42	126.07	111.30
1	A	127	ASP	CA-CB-CG	6.28	118.88	112.60
1	E	292	THR	CA-CB-OG1	-6.24	100.25	109.60
1	D	42	THR	N-CA-CB	-6.09	100.47	110.52
1	E	362	GLU	CB-CG-CD	5.98	122.77	112.60
1	H	148	ARG	CG-CD-NE	5.73	124.60	112.00
1	G	127	ASP	CA-CB-CG	5.67	118.27	112.60
1	D	342	ASP	CA-CB-CG	5.67	118.27	112.60
1	F	319	THR	CA-CB-OG1	-5.65	101.12	109.60
1	B	91	HIS	CA-CB-CG	5.63	119.43	113.80
1	E	288	TYR	CA-C-O	-5.62	114.47	120.42
1	F	165	ASP	CA-CB-CG	5.51	118.11	112.60
1	E	91	HIS	CA-CB-CG	5.50	119.30	113.80
1	A	113	THR	CB-CA-C	5.39	118.43	109.38
1	C	91	HIS	CA-CB-CG	5.21	119.02	113.80
1	H	295	ARG	CB-CG-CD	5.21	123.28	111.30
1	H	150	VAL	N-CA-C	-5.19	105.65	110.53

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	250	LEU	CA-C-N	5.18	127.22	120.28
1	E	250	LEU	C-N-CA	5.18	127.22	120.28
1	F	362	GLU	CB-CG-CD	5.15	121.35	112.60
1	G	345	ARG	NE-CZ-NH2	-5.06	114.64	119.20
1	C	42	THR	CA-CB-OG1	-5.05	102.02	109.60
1	C	361	ASP	CA-C-O	-5.04	115.53	120.82
1	G	91	HIS	CA-CB-CG	5.03	118.83	113.80
1	G	258	THR	CA-CB-OG1	-5.02	102.07	109.60
1	D	188	HIS	CA-C-O	-5.01	115.03	120.09
1	G	338	ASP	CA-CB-CG	5.01	117.61	112.60

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	20	TYR	Sidechain

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	2928	0	2752	37	0
1	B	2928	0	2752	28	0
1	C	2928	0	2752	18	0
1	D	2928	0	2752	16	0
1	E	2928	0	2752	24	0
1	F	2928	0	2752	29	0
1	G	2928	0	2752	25	0
1	H	2928	0	2752	26	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
2	E	1	0	0	0	0
2	F	1	0	0	0	0
2	G	1	0	0	0	0
2	H	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	A	177	0	0	9	2
3	B	181	0	0	5	2
3	C	182	0	0	4	0
3	D	171	0	0	6	0
3	E	105	0	0	5	0
3	F	95	0	0	5	0
3	G	96	0	0	6	0
3	H	103	0	0	6	0
All	All	24542	0	22016	189	2

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 4.

All (189) 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:G:42:THR:HG22	3:G:526:HOH:O	1.27	1.31
1:A:42:THR:HG22	3:A:585:HOH:O	1.29	1.25
1:C:42:THR:HG22	3:C:534:HOH:O	1.44	1.16
1:A:257:HIS:HD2	3:A:504:HOH:O	1.33	1.09
1:E:42:THR:HG22	3:E:523:HOH:O	1.54	1.08
1:H:42:THR:HG22	3:H:541:HOH:O	1.54	1.07
1:A:257:HIS:CD2	3:A:504:HOH:O	2.09	1.03
1:B:109:THR:HG22	1:B:148:ARG:HH12	1.19	1.03
1:A:368:LYS:O	1:A:372:MET:HG2	1.57	1.01
1:B:230:GLU:OE1	3:B:501:HOH:O	1.79	1.00
1:E:230:GLU:OE1	3:E:501:HOH:O	1.78	0.99
1:F:230:GLU:OE1	3:F:501:HOH:O	1.82	0.97
1:A:337:GLU:HB2	1:A:339:LEU:HD11	1.48	0.94
1:C:230:GLU:OE1	3:C:501:HOH:O	1.86	0.94
1:A:230:GLU:OE1	3:A:501:HOH:O	1.85	0.94
1:D:230:GLU:OE1	3:D:501:HOH:O	1.84	0.93
1:F:54:VAL:HG21	1:F:207:PHE:HZ	1.36	0.91
1:E:110:ALA:HB1	1:E:331:ASP:OD1	1.73	0.89
1:H:230:GLU:OE1	3:H:501:HOH:O	1.90	0.88
1:D:42:THR:HG23	3:D:552:HOH:O	1.74	0.86
1:A:54:VAL:HG21	1:A:207:PHE:HZ	1.42	0.84
1:A:71:MET:HA	1:A:71:MET:HE2	1.58	0.84
1:E:291:TYR:CZ	1:E:295:ARG:HD2	2.13	0.84
1:F:42:THR:HG23	3:F:527:HOH:O	1.78	0.82
1:G:230:GLU:OE1	3:G:501:HOH:O	1.95	0.82

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:109:THR:HG22	1:B:148:ARG:NH1	1.96	0.78
1:A:256:PHE:HB2	3:A:504:HOH:O	1.82	0.78
1:B:188:HIS:NE2	3:B:502:HOH:O	2.10	0.77
1:C:109:THR:CG2	1:C:148:ARG:HH12	1.98	0.76
1:B:54:VAL:HG21	1:B:207:PHE:HZ	1.50	0.75
1:F:54:VAL:HG21	1:F:207:PHE:CZ	2.21	0.75
1:G:55:GLU:OE1	1:G:61:LYS:HE2	1.88	0.73
1:E:331:ASP:OD1	3:E:502:HOH:O	2.05	0.73
1:A:86:LYS:NZ	3:A:503:HOH:O	2.20	0.73
1:C:109:THR:HG23	1:C:148:ARG:HH12	1.52	0.72
1:H:54:VAL:HG21	1:H:207:PHE:HZ	1.54	0.72
1:E:291:TYR:OH	1:E:295:ARG:HD2	1.90	0.72
1:A:54:VAL:HG21	1:A:207:PHE:CZ	2.24	0.72
1:D:291:TYR:CZ	1:D:295:ARG:HD2	2.25	0.71
1:D:188:HIS:NE2	3:D:502:HOH:O	2.11	0.71
1:A:337:GLU:CB	1:A:339:LEU:HD11	2.19	0.70
1:E:188:HIS:NE2	3:E:503:HOH:O	2.09	0.70
1:E:168:ALA:HB1	1:E:171:GLU:HG3	1.72	0.69
1:H:42:THR:CG2	3:H:541:HOH:O	2.25	0.69
1:A:71:MET:HE2	1:A:74:ILE:HD12	1.73	0.69
1:C:188:HIS:NE2	3:C:502:HOH:O	2.15	0.69
1:G:110:ALA:HB1	1:G:331:ASP:OD1	1.92	0.69
1:G:188:HIS:NE2	3:G:502:HOH:O	2.04	0.68
1:G:251:ARG:HD3	1:G:255:ASP:OD2	1.96	0.67
1:F:238:MET:HE2	1:F:252:ILE:HD13	1.76	0.66
1:H:342:ASP:OD2	1:H:344:ASN:HB2	1.95	0.66
1:A:337:GLU:HB2	1:A:339:LEU:CD1	2.25	0.66
1:D:198:LEU:C	1:D:198:LEU:HD23	2.22	0.65
1:B:54:VAL:HG21	1:B:207:PHE:CZ	2.31	0.64
1:G:331:ASP:OD1	3:G:503:HOH:O	2.14	0.64
1:A:144:ASP:OD2	1:B:94:THR:HG23	1.97	0.64
1:B:109:THR:CG2	1:B:148:ARG:HH22	2.11	0.64
1:A:71:MET:CE	1:A:74:ILE:HD12	2.27	0.64
1:E:144:ASP:OD2	1:F:94:THR:HG23	1.98	0.63
1:C:144:ASP:OD2	1:D:94:THR:HG23	1.99	0.63
1:H:54:VAL:HG21	1:H:207:PHE:CZ	2.33	0.63
1:F:188:HIS:NE2	3:F:502:HOH:O	2.09	0.62
1:A:66:ILE:HD11	1:A:71:MET:CE	2.29	0.62
1:H:238:MET:HE2	1:H:252:ILE:HD13	1.79	0.62
1:E:335:PRO:O	1:E:336:SER:O	2.19	0.61
1:B:230:GLU:CD	3:B:501:HOH:O	2.37	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:40:MET:HG3	1:B:281:PHE:CE1	2.36	0.60
1:E:109:THR:HG22	1:E:148:ARG:HE	1.66	0.59
1:D:167:ARG:NE	3:D:504:HOH:O	2.34	0.59
1:F:362:GLU:CD	1:F:362:GLU:H	2.10	0.59
1:G:94:THR:HG23	1:H:144:ASP:OD2	2.02	0.59
1:G:343:ARG:O	1:G:343:ARG:HD3	2.02	0.59
1:G:54:VAL:HG21	1:G:207:PHE:HZ	1.67	0.59
1:F:40:MET:HB3	1:F:41:PRO:HD2	1.85	0.58
1:E:109:THR:CG2	1:E:148:ARG:HH21	2.17	0.58
1:F:109:THR:HG23	1:F:148:ARG:HH12	1.69	0.57
1:B:368:LYS:HG2	1:B:372:MET:CE	2.34	0.57
1:G:54:VAL:HG21	1:G:207:PHE:CZ	2.40	0.57
1:G:295:ARG:NH2	1:G:300:GLY:O	2.25	0.57
1:A:71:MET:HE2	1:A:71:MET:CA	2.32	0.56
1:D:167:ARG:CZ	3:D:504:HOH:O	2.53	0.56
1:H:82:ASN:ND2	1:H:336:SER:OG	2.38	0.56
1:D:40:MET:HB3	1:D:41:PRO:HD2	1.86	0.56
1:G:42:THR:CG2	3:G:526:HOH:O	2.08	0.56
1:F:109:THR:CG2	1:F:148:ARG:HH12	2.19	0.56
1:H:342:ASP:OD2	1:H:344:ASN:CB	2.53	0.55
1:F:152:ILE:HG22	1:F:153:ASN:N	2.21	0.55
1:A:66:ILE:HD11	1:A:71:MET:HE1	1.87	0.55
1:E:362:GLU:H	1:E:362:GLU:CD	2.14	0.55
1:H:109:THR:CG2	1:H:148:ARG:HH21	2.19	0.55
1:H:40:MET:HB3	1:H:41:PRO:HD2	1.89	0.54
1:H:141:TYR:OH	3:H:502:HOH:O	2.03	0.54
1:A:249:ALA:HA	1:A:252:ILE:HD12	1.90	0.54
1:E:110:ALA:HB1	1:E:331:ASP:CG	2.33	0.53
1:E:54:VAL:HG21	1:E:207:PHE:CZ	2.44	0.53
1:C:54:VAL:HG21	1:C:207:PHE:CZ	2.43	0.53
1:C:198:LEU:C	1:C:198:LEU:HD23	2.34	0.53
1:B:368:LYS:HG2	1:B:372:MET:HE2	1.90	0.53
1:D:291:TYR:OH	1:D:295:ARG:HD2	2.08	0.52
1:F:54:VAL:CG2	1:F:207:PHE:HZ	2.17	0.52
1:C:168:ALA:HB1	1:C:171:GLU:HG3	1.92	0.52
1:B:54:VAL:CG2	1:B:207:PHE:HZ	2.22	0.52
1:H:155:ARG:HD3	1:H:339:LEU:HD21	1.91	0.51
1:D:250:LEU:O	1:D:254:LEU:HD23	2.11	0.51
1:F:152:ILE:CG2	1:F:153:ASN:N	2.74	0.51
1:H:109:THR:HG23	1:H:148:ARG:HH21	1.76	0.51
1:A:14:PRO:HB2	1:A:15:LYS:HE3	1.93	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:198:LEU:C	1:E:198:LEU:HD23	2.36	0.51
1:E:254:LEU:HD21	1:F:368:LYS:HD3	1.93	0.50
1:B:362:GLU:H	1:B:362:GLU:CD	2.20	0.50
1:A:339:LEU:HD12	1:A:339:LEU:N	2.26	0.50
1:A:66:ILE:HD11	1:A:71:MET:HE3	1.94	0.49
1:H:188:HIS:NE2	3:H:503:HOH:O	2.06	0.49
1:C:109:THR:HG23	1:C:148:ARG:NH1	2.25	0.49
1:G:257:HIS:ND1	3:G:501:HOH:O	2.35	0.49
1:C:42:THR:CG2	3:C:534:HOH:O	2.24	0.49
1:D:40:MET:HB3	1:D:41:PRO:CD	2.42	0.49
1:A:71:MET:HA	1:A:71:MET:CE	2.37	0.49
1:A:372:MET:HE1	1:B:250:LEU:CD2	2.43	0.49
1:E:54:VAL:HG21	1:E:207:PHE:HZ	1.78	0.49
1:F:342:ASP:N	1:F:342:ASP:OD1	2.45	0.49
1:F:42:THR:CG2	3:F:527:HOH:O	2.47	0.49
1:B:42:THR:HG23	3:B:544:HOH:O	2.12	0.48
1:F:233:MET:HE1	1:F:242:TYR:HE2	1.77	0.48
1:C:250:LEU:HD21	1:D:372:MET:HE3	1.95	0.48
1:G:54:VAL:CG2	1:G:207:PHE:HZ	2.27	0.48
1:A:337:GLU:CB	1:A:339:LEU:CD1	2.89	0.47
1:D:230:GLU:CD	3:D:501:HOH:O	2.44	0.47
1:G:19:TYR:O	1:G:323:ALA:HA	2.14	0.47
1:A:216:TRP:CH2	1:B:356:GLY:HA3	2.50	0.47
1:A:42:THR:CG2	3:A:585:HOH:O	2.12	0.47
1:C:109:THR:HG22	1:C:148:ARG:HH12	1.75	0.47
1:B:155:ARG:HB3	1:B:156:PRO:HD3	1.97	0.47
1:H:291:TYR:CZ	1:H:295:ARG:HD2	2.49	0.47
1:C:40:MET:HB3	1:C:41:PRO:HD2	1.98	0.46
1:F:109:THR:CG2	1:F:148:ARG:HH22	2.27	0.46
1:E:42:THR:CG2	3:E:523:HOH:O	2.35	0.46
1:F:198:LEU:C	1:F:198:LEU:HD23	2.40	0.46
1:H:342:ASP:CG	1:H:344:ASN:HB2	2.41	0.46
1:E:40:MET:HG3	1:E:281:PHE:CE1	2.51	0.46
1:F:37:GLY:HA3	1:F:49:ALA:O	2.15	0.46
1:B:198:LEU:C	1:B:198:LEU:HD23	2.41	0.45
1:F:230:GLU:CD	3:F:501:HOH:O	2.47	0.45
1:G:362:GLU:H	1:G:362:GLU:CD	2.24	0.45
1:D:233:MET:HE1	1:D:242:TYR:HE2	1.81	0.45
1:A:188:HIS:NE2	3:A:502:HOH:O	2.08	0.45
1:A:372:MET:HE1	1:B:250:LEU:HD21	1.99	0.45
1:A:251:ARG:HD3	1:A:255:ASP:OD2	2.17	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:110:ALA:HB1	1:G:331:ASP:CG	2.41	0.44
1:B:40:MET:SD	1:B:233:MET:HE2	2.56	0.44
1:C:54:VAL:HG21	1:C:207:PHE:HZ	1.82	0.44
1:A:40:MET:HB3	1:A:41:PRO:CD	2.48	0.43
1:B:40:MET:HG3	1:B:281:PHE:CD1	2.54	0.43
1:H:199:GLU:CD	3:H:526:HOH:O	2.61	0.43
1:F:342:ASP:OD2	1:F:343:ARG:NH1	2.51	0.43
1:A:252:ILE:HG23	3:A:504:HOH:O	2.19	0.43
1:H:291:TYR:OH	1:H:295:ARG:HD2	2.19	0.43
1:A:198:LEU:C	1:A:198:LEU:HD23	2.44	0.43
1:G:144:ASP:OD2	1:H:94:THR:HG23	2.18	0.43
1:B:338:ASP:O	1:B:339:LEU:HB2	2.18	0.42
1:C:238:MET:HE2	1:C:252:ILE:HD13	2.01	0.42
1:E:238:MET:HE2	1:E:252:ILE:HD13	2.01	0.42
1:A:192:GLY:N	1:A:193:PRO:CD	2.82	0.42
1:C:251:ARG:HD3	1:C:255:ASP:OD2	2.20	0.42
1:F:362:GLU:CD	1:F:362:GLU:N	2.75	0.42
1:B:339:LEU:HD12	1:B:339:LEU:HA	1.93	0.42
1:F:287:ARG:HD3	1:H:288:TYR:OH	2.20	0.41
1:H:40:MET:HB3	1:H:41:PRO:CD	2.49	0.41
1:D:198:LEU:C	1:D:198:LEU:CD2	2.91	0.41
1:F:247:GLN:HG3	1:F:248:LEU:HD12	2.02	0.41
1:B:325:SER:HB2	1:B:326:PRO:HD2	2.02	0.41
1:H:290:ILE:HG22	1:H:304:MET:HE1	2.03	0.41
1:G:291:TYR:CZ	1:E:288:TYR:HB3	2.55	0.41
1:B:40:MET:HB3	1:B:41:PRO:HD2	2.02	0.41
1:H:5:LYS:HB3	1:H:5:LYS:HE3	1.76	0.41
1:A:37:GLY:HA3	1:A:49:ALA:O	2.21	0.41
1:G:54:VAL:CG2	1:G:207:PHE:CZ	3.03	0.41
1:F:101:VAL:HG21	1:F:141:TYR:CE2	2.56	0.41
1:F:152:ILE:HG22	1:F:153:ASN:O	2.21	0.41
1:B:342:ASP:O	3:B:503:HOH:O	2.22	0.41
1:B:368:LYS:CG	1:B:372:MET:HE2	2.51	0.41
1:G:288:TYR:HB3	1:E:291:TYR:CZ	2.56	0.41
1:C:19:TYR:O	1:C:323:ALA:HA	2.20	0.41
1:F:19:TYR:O	1:F:323:ALA:HA	2.21	0.41
1:E:105:ALA:O	1:E:109:THR:HG23	2.21	0.40
1:G:238:MET:HE2	1:G:252:ILE:HD13	2.03	0.40
1:G:356:GLY:HA3	1:H:216:TRP:CH2	2.57	0.40
1:G:291:TYR:CE1	1:G:302:LEU:HD23	2.57	0.40
1:A:108:TRP:CE3	1:A:148:ARG:HD2	2.57	0.40

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

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:A:665:HOH:O	3:B:654:HOH:O[2_555]	2.03	0.17
3:A:518:HOH:O	3:B:551:HOH:O[2_545]	2.06	0.14

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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	370/372 (100%)	358 (97%)	11 (3%)	1 (0%)	36	20
1	B	370/372 (100%)	357 (96%)	9 (2%)	4 (1%)	11	2
1	C	370/372 (100%)	360 (97%)	9 (2%)	1 (0%)	36	20
1	D	370/372 (100%)	359 (97%)	10 (3%)	1 (0%)	36	20
1	E	370/372 (100%)	360 (97%)	8 (2%)	2 (0%)	24	10
1	F	370/372 (100%)	359 (97%)	10 (3%)	1 (0%)	36	20
1	G	370/372 (100%)	357 (96%)	11 (3%)	2 (0%)	24	10
1	H	370/372 (100%)	356 (96%)	10 (3%)	4 (1%)	11	2
All	All	2960/2976 (100%)	2866 (97%)	78 (3%)	16 (0%)	24	10

All (16) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	G	336	SER
1	E	336	SER
1	B	339	LEU
1	H	341	PRO
1	H	338	ASP
1	H	344	ASN
1	G	66	ILE
1	B	340	ALA

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Mol	Chain	Res	Type
1	H	66	ILE
1	A	66	ILE
1	D	66	ILE
1	B	66	ILE
1	B	337	GLU
1	C	66	ILE
1	E	66	ILE
1	F	66	ILE

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	301/301 (100%)	285 (95%)	16 (5%)	20	4
1	B	301/301 (100%)	294 (98%)	7 (2%)	44	20
1	C	301/301 (100%)	293 (97%)	8 (3%)	39	15
1	D	301/301 (100%)	292 (97%)	9 (3%)	36	12
1	E	301/301 (100%)	290 (96%)	11 (4%)	30	8
1	F	301/301 (100%)	293 (97%)	8 (3%)	39	15
1	G	301/301 (100%)	291 (97%)	10 (3%)	33	11
1	H	301/301 (100%)	292 (97%)	9 (3%)	36	12
All	All	2408/2408 (100%)	2330 (97%)	78 (3%)	34	11

All (78) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	5	LYS
1	A	15	LYS
1	A	40	MET
1	A	71	MET
1	A	86	LYS
1	A	103	MET
1	A	131	LYS

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Mol	Chain	Res	Type
1	A	169	LEU
1	A	176	GLU
1	A	230	GLU
1	A	244	GLN
1	A	247	GLN
1	A	295	ARG
1	A	368	LYS
1	A	369	GLN
1	A	372	MET
1	G	131	LYS
1	G	169	LEU
1	G	248	LEU
1	G	252	ILE
1	G	254	LEU
1	G	295	ARG
1	G	336	SER
1	G	337	GLU
1	G	338	ASP
1	G	362	GLU
1	C	32	ARG
1	C	40	MET
1	C	169	LEU
1	C	222	LYS
1	C	254	LEU
1	C	338	ASP
1	C	343	ARG
1	C	362	GLU
1	E	15	LYS
1	E	40	MET
1	E	131	LYS
1	E	169	LEU
1	E	254	LEU
1	E	295	ARG
1	E	316	ASP
1	E	337	GLU
1	E	343	ARG
1	E	362	GLU
1	E	369	GLN
1	D	42	THR
1	D	131	LYS
1	D	169	LEU
1	D	175	GLN

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Mol	Chain	Res	Type
1	D	252	ILE
1	D	254	LEU
1	D	295	ARG
1	D	337	GLU
1	D	339	LEU
1	F	12	ILE
1	F	42	THR
1	F	169	LEU
1	F	247	GLN
1	F	254	LEU
1	F	295	ARG
1	F	342	ASP
1	F	362	GLU
1	B	109	THR
1	B	169	LEU
1	B	222	LYS
1	B	248	LEU
1	B	254	LEU
1	B	336	SER
1	B	362	GLU
1	H	15	LYS
1	H	148	ARG
1	H	152	ILE
1	H	169	LEU
1	H	247	GLN
1	H	254	LEU
1	H	295	ARG
1	H	337	GLU
1	H	362	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (14) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	344	ASN
1	G	344	ASN
1	E	170	ASN
1	E	175	GLN
1	D	170	ASN
1	F	170	ASN
1	F	247	GLN
1	B	170	ASN
1	B	175	GLN

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Mol	Chain	Res	Type
1	B	244	GLN
1	H	82	ASN
1	H	153	ASN
1	H	170	ASN
1	H	344	ASN

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 8 ligands modelled in this entry, 8 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 [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	372/372 (100%)	0.75	27 (7%)	21 22	6, 14, 30, 50	0
1	B	372/372 (100%)	0.26	37 (9%)	13 13	4, 10, 40, 84	0
1	C	372/372 (100%)	0.03	24 (6%)	25 27	3, 8, 32, 69	0
1	D	372/372 (100%)	0.36	25 (6%)	24 25	5, 12, 30, 60	0
1	E	372/372 (100%)	0.10	27 (7%)	21 22	3, 9, 30, 71	0
1	F	372/372 (100%)	0.30	27 (7%)	21 22	5, 12, 31, 49	0
1	G	372/372 (100%)	0.89	54 (14%)	6 6	6, 15, 31, 61	0
1	H	372/372 (100%)	0.35	40 (10%)	11 11	4, 10, 41, 84	0
All	All	2976/2976 (100%)	0.38	261 (8%)	15 16	3, 11, 34, 84	0

All (261) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	H	152	ILE	12.4
1	B	339	LEU	8.7
1	H	151	THR	8.1
1	H	248	LEU	7.7
1	H	341	PRO	7.3
1	H	339	LEU	7.0
1	B	373	GLY	6.9
1	F	248	LEU	6.7
1	F	339	LEU	6.6
1	E	373	GLY	6.2
1	F	373	GLY	6.2
1	H	342	ASP	6.2
1	G	338	ASP	6.1
1	H	343	ARG	6.1
1	D	339	LEU	5.9
1	G	248	LEU	5.8

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Mol	Chain	Res	Type	RSRZ
1	A	373	GLY	5.8
1	E	340	ALA	5.7
1	H	344	ASN	5.7
1	H	340	ALA	5.5
1	D	336	SER	5.5
1	E	337	GLU	5.4
1	H	373	GLY	5.3
1	A	254	LEU	5.3
1	G	337	GLU	5.3
1	H	338	ASP	5.3
1	B	338	ASP	5.2
1	E	339	LEU	5.2
1	B	248	LEU	5.2
1	G	339	LEU	5.1
1	C	338	ASP	5.1
1	D	373	GLY	4.9
1	C	339	LEU	4.9
1	F	251	ARG	4.8
1	G	373	GLY	4.8
1	B	152	ILE	4.7
1	E	343	ARG	4.7
1	B	336	SER	4.7
1	H	153	ASN	4.6
1	H	336	SER	4.6
1	E	338	ASP	4.6
1	F	152	ILE	4.6
1	C	337	GLU	4.5
1	C	340	ALA	4.5
1	C	343	ARG	4.5
1	F	336	SER	4.4
1	G	336	SER	4.4
1	B	369	GLN	4.3
1	H	337	GLU	4.3
1	E	243	ASN	4.2
1	H	243	ASN	4.2
1	H	366	HIS	4.2
1	F	12	ILE	4.1
1	B	340	ALA	3.9
1	C	362	GLU	3.9
1	B	337	GLU	3.9
1	B	371	PHE	3.9
1	C	373	GLY	3.8

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Mol	Chain	Res	Type	RSRZ
1	D	254	LEU	3.8
1	C	366	HIS	3.8
1	G	372	MET	3.8
1	B	342	ASP	3.7
1	E	331	ASP	3.7
1	B	243	ASN	3.7
1	B	366	HIS	3.7
1	G	316	ASP	3.6
1	F	338	ASP	3.6
1	H	250	LEU	3.6
1	B	247	GLN	3.6
1	C	251	ARG	3.6
1	A	248	LEU	3.6
1	H	247	GLN	3.5
1	E	256	PHE	3.5
1	H	371	PHE	3.5
1	H	256	PHE	3.5
1	H	254	LEU	3.4
1	G	366	HIS	3.4
1	H	363	ALA	3.4
1	D	243	ASN	3.4
1	D	337	GLU	3.3
1	E	251	ARG	3.3
1	G	317	ALA	3.3
1	G	362	GLU	3.3
1	D	366	HIS	3.3
1	F	366	HIS	3.3
1	B	246	ALA	3.3
1	B	352	TYR	3.2
1	H	369	GLN	3.2
1	A	245	PRO	3.2
1	E	344	ASN	3.2
1	G	314	THR	3.2
1	G	343	ARG	3.2
1	B	254	LEU	3.2
1	A	256	PHE	3.2
1	C	256	PHE	3.2
1	A	343	ARG	3.2
1	H	251	ARG	3.1
1	G	335	PRO	3.1
1	E	316	ASP	3.1
1	C	248	LEU	3.1

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Mol	Chain	Res	Type	RSRZ
1	A	369	GLN	3.1
1	G	292	THR	3.1
1	A	251	ARG	3.1
1	B	362	GLU	3.1
1	A	372	MET	3.1
1	D	251	ARG	3.1
1	E	362	GLU	3.1
1	E	341	PRO	3.0
1	H	245	PRO	3.0
1	H	362	GLU	3.0
1	E	366	HIS	3.0
1	G	12	ILE	3.0
1	H	246	ALA	3.0
1	G	29	ASP	3.0
1	F	342	ASP	3.0
1	G	110	ALA	3.0
1	G	254	LEU	3.0
1	F	337	GLU	3.0
1	D	342	ASP	2.9
1	G	291	TYR	2.9
1	A	243	ASN	2.9
1	F	243	ASN	2.9
1	A	247	GLN	2.9
1	G	371	PHE	2.9
1	F	340	ALA	2.9
1	H	249	ALA	2.9
1	C	341	PRO	2.9
1	F	369	GLN	2.9
1	G	240	THR	2.9
1	F	151	THR	2.9
1	F	246	ALA	2.9
1	C	254	LEU	2.9
1	H	150	VAL	2.8
1	B	367	TRP	2.8
1	H	367	TRP	2.8
1	G	109	THR	2.8
1	F	254	LEU	2.8
1	C	344	ASN	2.8
1	G	247	GLN	2.8
1	B	251	ARG	2.8
1	B	250	LEU	2.8
1	F	109	THR	2.8

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Mol	Chain	Res	Type	RSRZ
1	G	299	ALA	2.7
1	E	244	GLN	2.7
1	B	256	PHE	2.7
1	A	169	LEU	2.7
1	A	250	LEU	2.7
1	E	254	LEU	2.7
1	E	336	SER	2.7
1	A	246	ALA	2.7
1	H	368	LYS	2.7
1	E	335	PRO	2.7
1	C	244	GLN	2.7
1	C	247	GLN	2.7
1	A	366	HIS	2.7
1	E	369	GLN	2.7
1	D	369	GLN	2.7
1	D	338	ASP	2.7
1	G	365	ALA	2.7
1	F	362	GLU	2.6
1	B	370	GLU	2.6
1	C	335	PRO	2.6
1	A	371	PHE	2.6
1	E	247	GLN	2.6
1	B	365	ALA	2.6
1	G	222	LYS	2.6
1	A	110	ALA	2.6
1	B	363	ALA	2.6
1	G	369	GLN	2.5
1	E	248	LEU	2.5
1	E	372	MET	2.5
1	B	151	THR	2.5
1	C	369	GLN	2.5
1	F	244	GLN	2.5
1	E	246	ALA	2.5
1	D	335	PRO	2.5
1	B	150	VAL	2.5
1	B	249	ALA	2.5
1	H	372	MET	2.5
1	A	176	GLU	2.5
1	G	155	ARG	2.5
1	G	251	ARG	2.5
1	G	315	ARG	2.5
1	F	372	MET	2.5

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Mol	Chain	Res	Type	RSRZ
1	C	371	PHE	2.4
1	D	256	PHE	2.4
1	G	363	ALA	2.4
1	G	270	VAL	2.4
1	A	370	GLU	2.4
1	H	359	ASN	2.4
1	G	111	GLY	2.4
1	B	245	PRO	2.4
1	G	200	TRP	2.4
1	C	252	ILE	2.4
1	G	294	VAL	2.4
1	A	109	THR	2.3
1	G	221	ALA	2.3
1	G	331	ASP	2.3
1	D	29	ASP	2.3
1	D	340	ALA	2.3
1	D	368	LYS	2.3
1	C	372	MET	2.3
1	G	264	GLY	2.3
1	G	28	PRO	2.3
1	B	335	PRO	2.3
1	H	335	PRO	2.3
1	D	363	ALA	2.3
1	B	252	ILE	2.3
1	H	252	ILE	2.3
1	G	293	GLY	2.3
1	G	295	ARG	2.3
1	G	368	LYS	2.3
1	B	372	MET	2.3
1	D	244	GLN	2.2
1	B	255	ASP	2.2
1	D	360	VAL	2.2
1	A	240	THR	2.2
1	F	7	THR	2.2
1	C	243	ASN	2.2
1	D	246	ALA	2.2
1	H	365	ALA	2.2
1	E	250	LEU	2.2
1	F	11	GLY	2.2
1	G	203	TYR	2.2
1	F	335	PRO	2.2
1	B	240	THR	2.2

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Mol	Chain	Res	Type	RSRZ
1	H	364	ASN	2.2
1	G	332	VAL	2.2
1	G	175	GLN	2.2
1	A	291	TYR	2.2
1	A	368	LYS	2.2
1	B	341	PRO	2.2
1	E	332	VAL	2.1
1	G	201	ASN	2.1
1	A	316	ASP	2.1
1	B	239	MET	2.1
1	A	252	ILE	2.1
1	F	365	ALA	2.1
1	G	223	GLY	2.1
1	H	15	LYS	2.1
1	G	250	LEU	2.1
1	C	250	LEU	2.1
1	D	250	LEU	2.1
1	A	269	MET	2.1
1	A	295	ARG	2.1
1	D	341	PRO	2.1
1	E	368	LYS	2.1
1	D	372	MET	2.1
1	F	250	LEU	2.1
1	G	252	ILE	2.1
1	G	370	GLU	2.0
1	G	233	MET	2.0
1	B	357	ARG	2.0
1	H	360	VAL	2.0
1	H	109	THR	2.0
1	G	268	ASN	2.0
1	F	4	GLY	2.0
1	G	290	ILE	2.0
1	D	371	PHE	2.0
1	C	365	ALA	2.0
1	D	255	ASP	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](#)

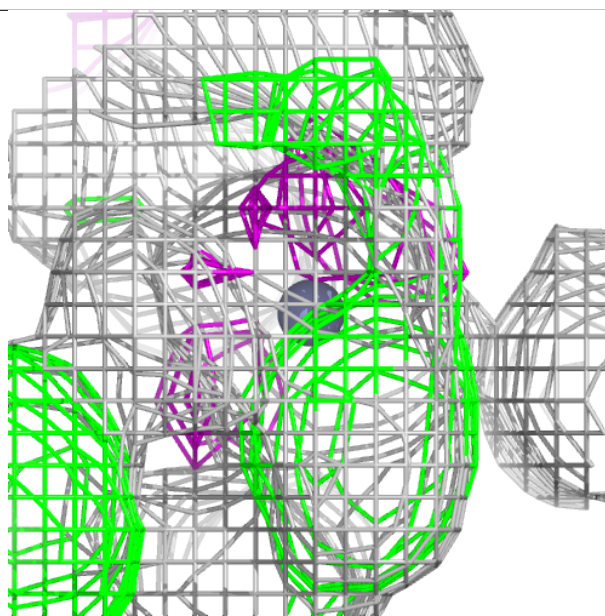
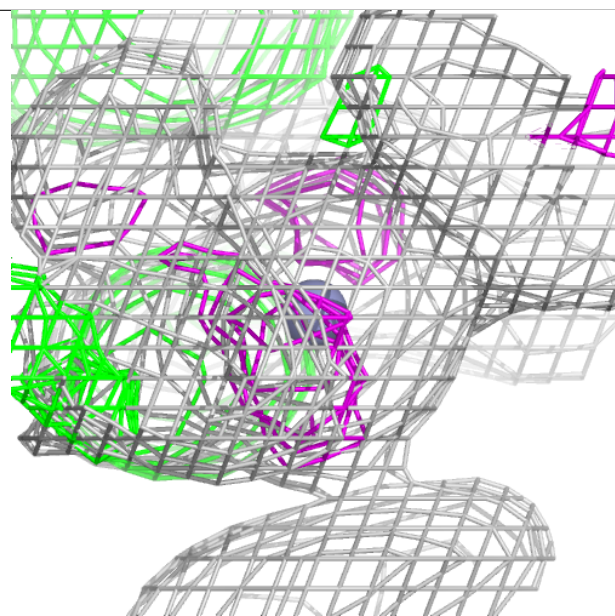
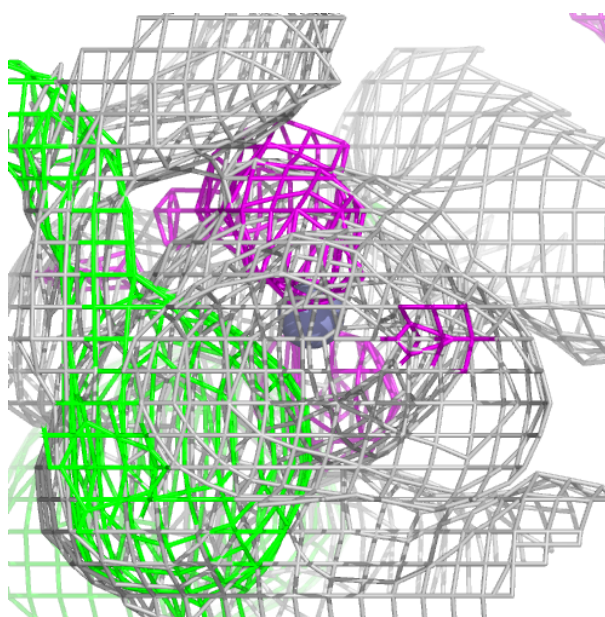
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	ZN	A	401	1/1	0.98	0.03	11,11,11,11	0
2	ZN	G	401	1/1	0.98	0.04	10,10,10,10	0
2	ZN	E	401	1/1	0.99	0.02	8,8,8,8	0
2	ZN	D	401	1/1	0.99	0.02	9,9,9,9	0
2	ZN	B	401	1/1	0.99	0.01	9,9,9,9	0
2	ZN	H	401	1/1	0.99	0.02	9,9,9,9	0
2	ZN	C	401	1/1	1.00	0.01	7,7,7,7	0
2	ZN	F	401	1/1	1.00	0.01	8,8,8,8	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.

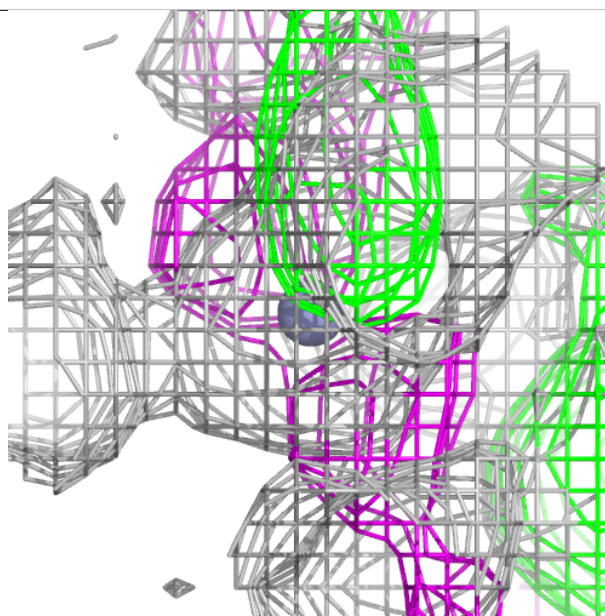
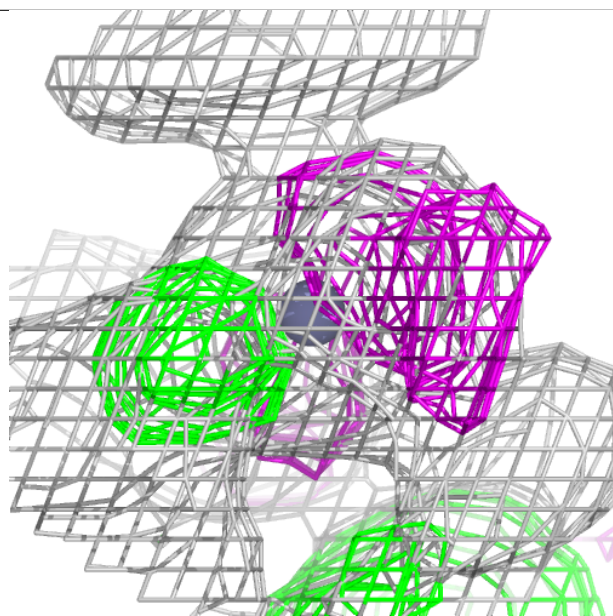
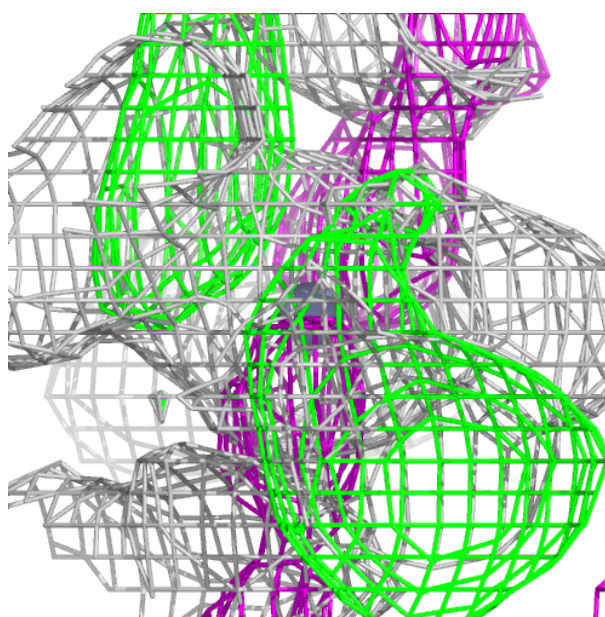
Electron density around ZN A 401:

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



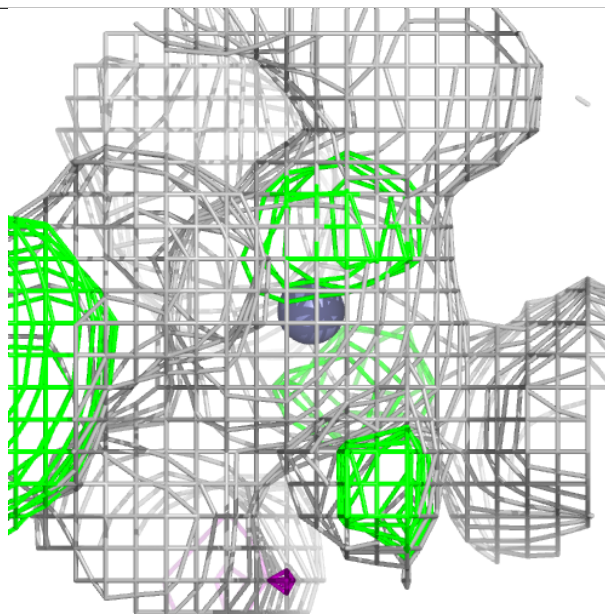
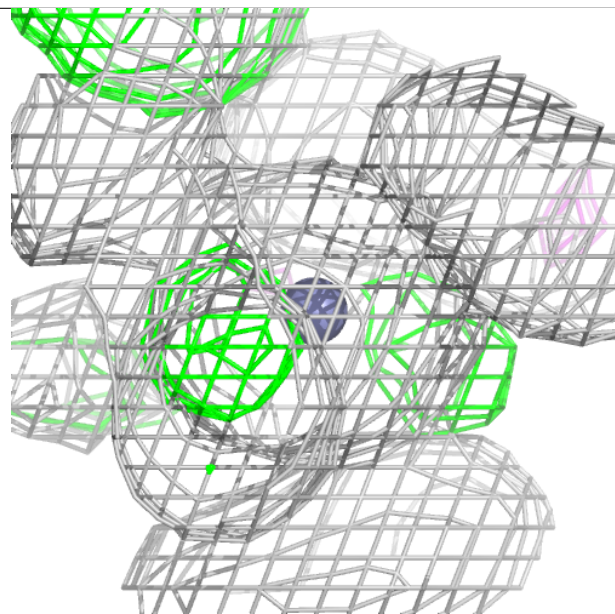
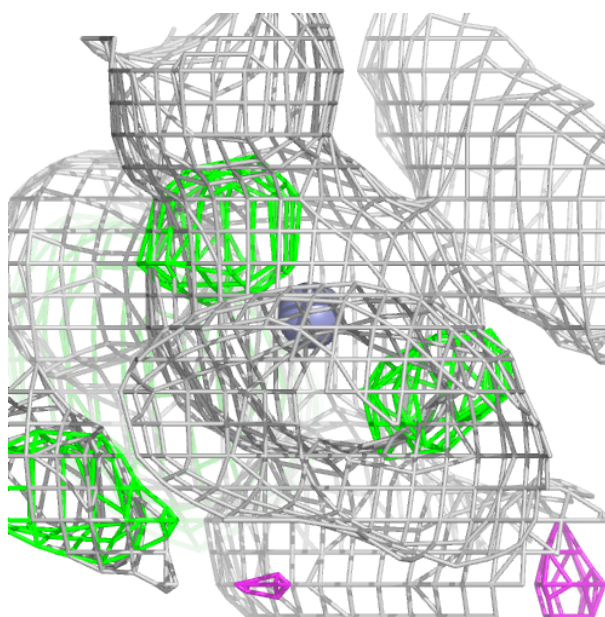
Electron density around ZN G 401:

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



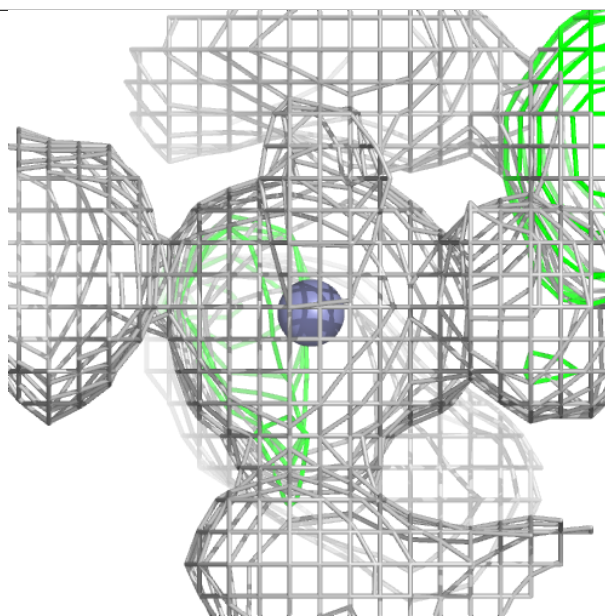
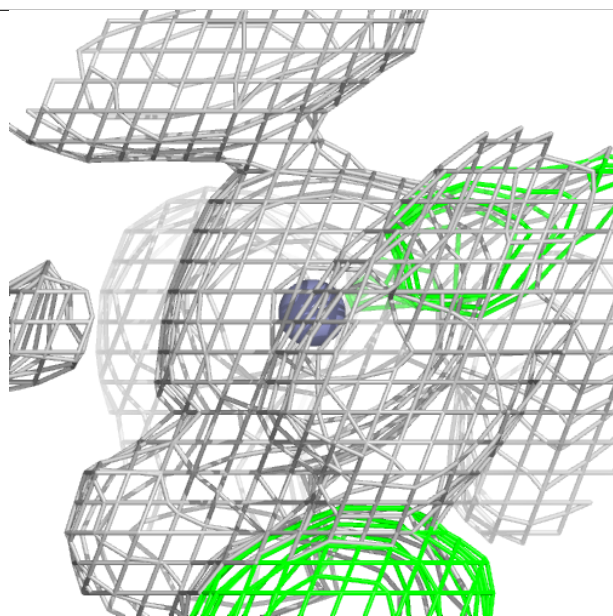
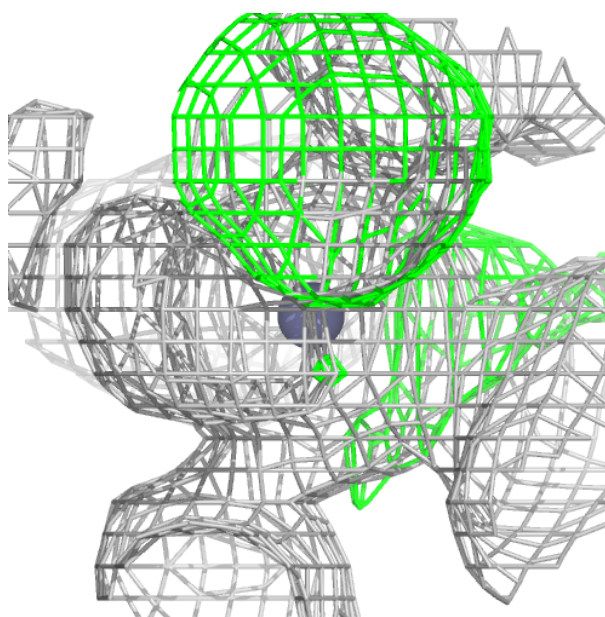
Electron density around ZN E 401:

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



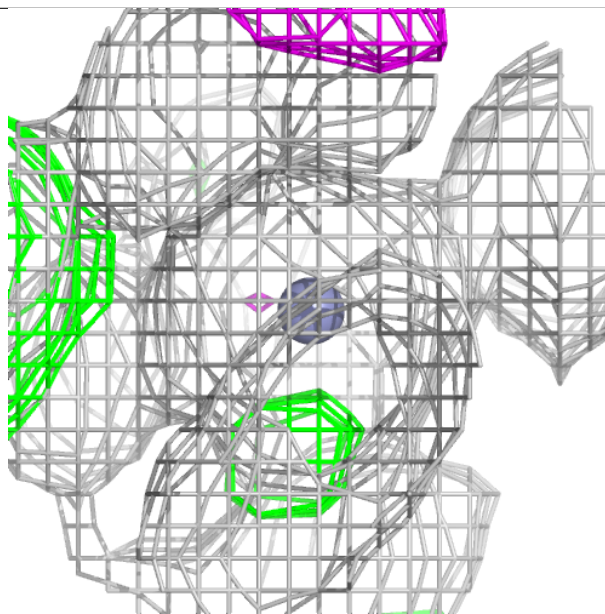
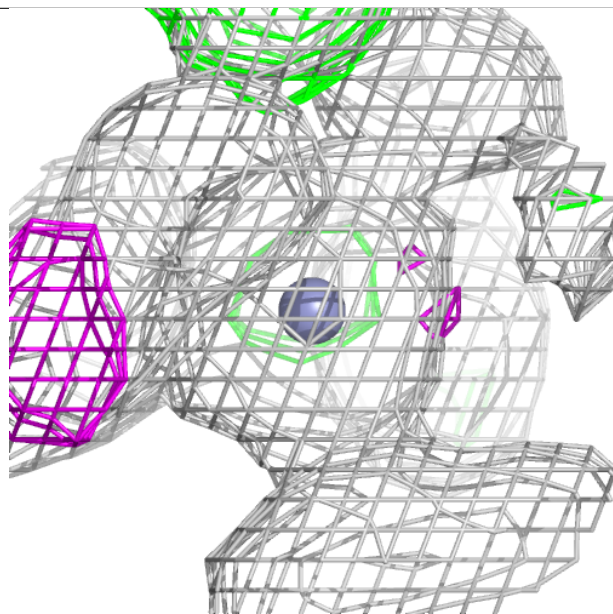
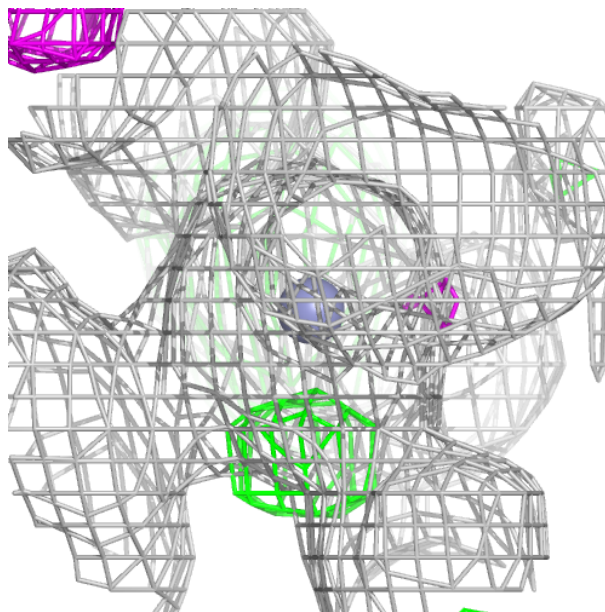
Electron density around ZN D 401:

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 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



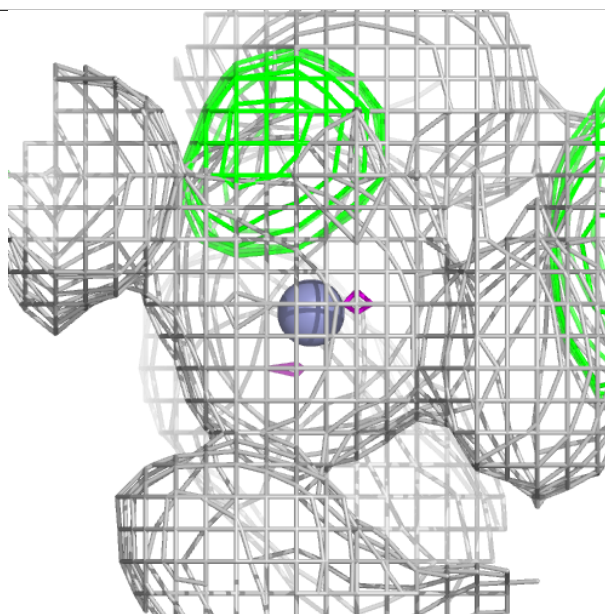
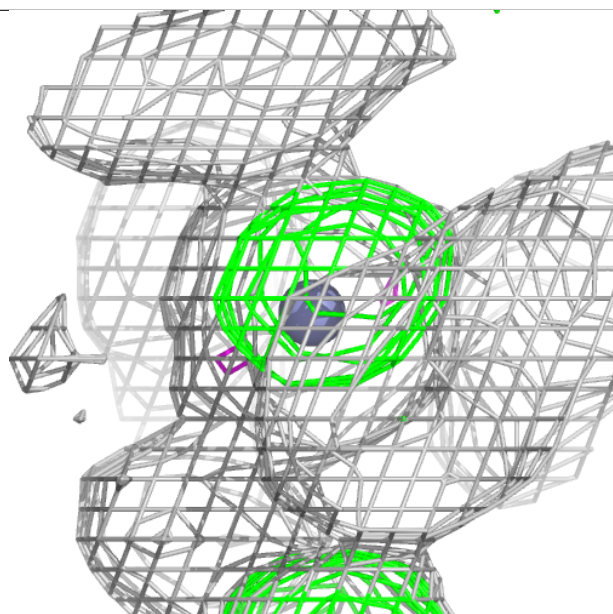
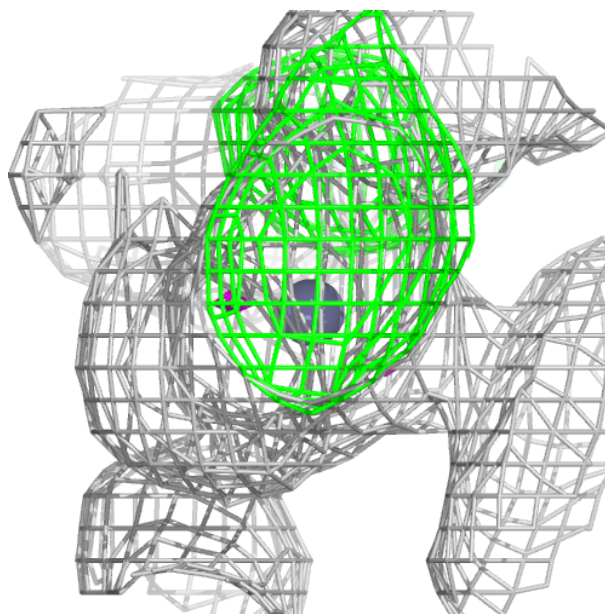
Electron density around ZN B 401:

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



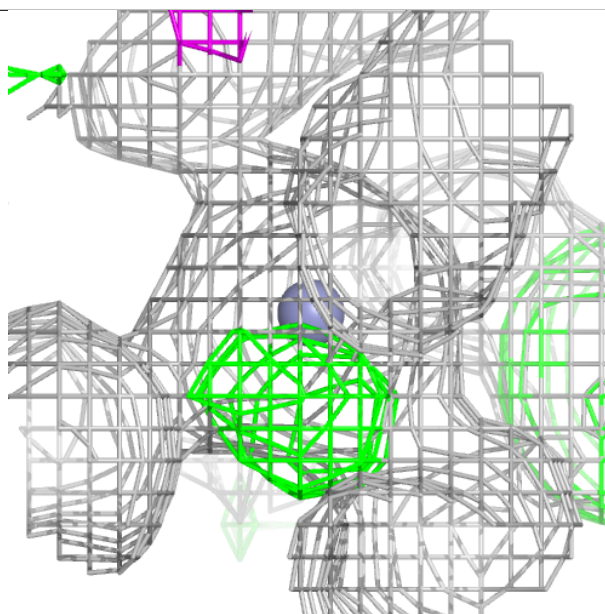
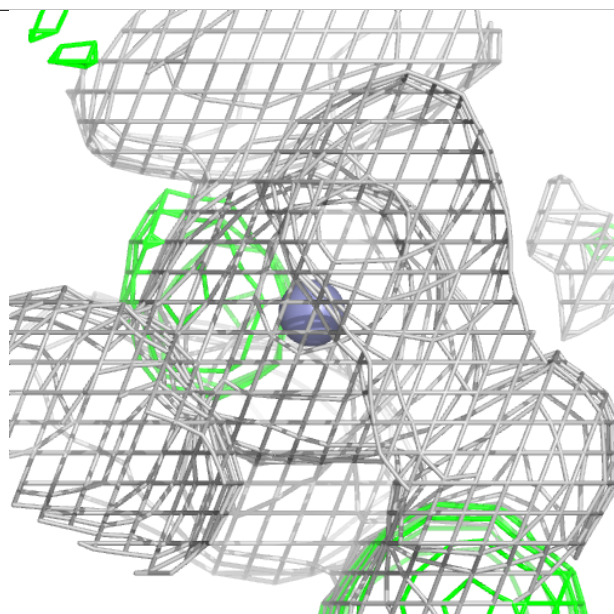
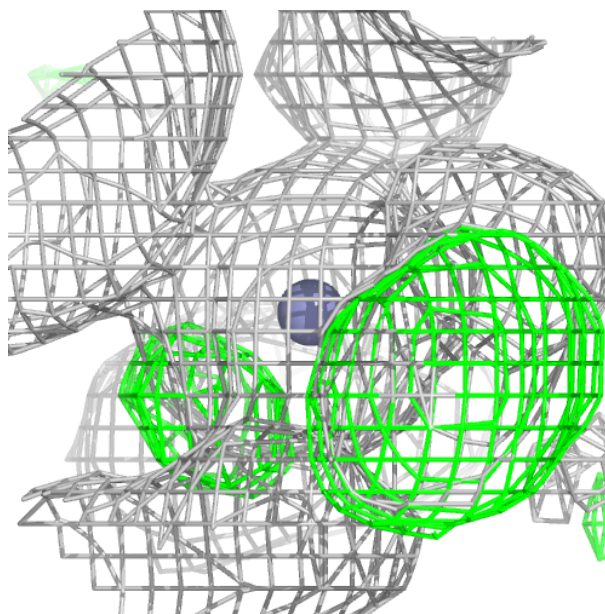
Electron density around ZN H 401:

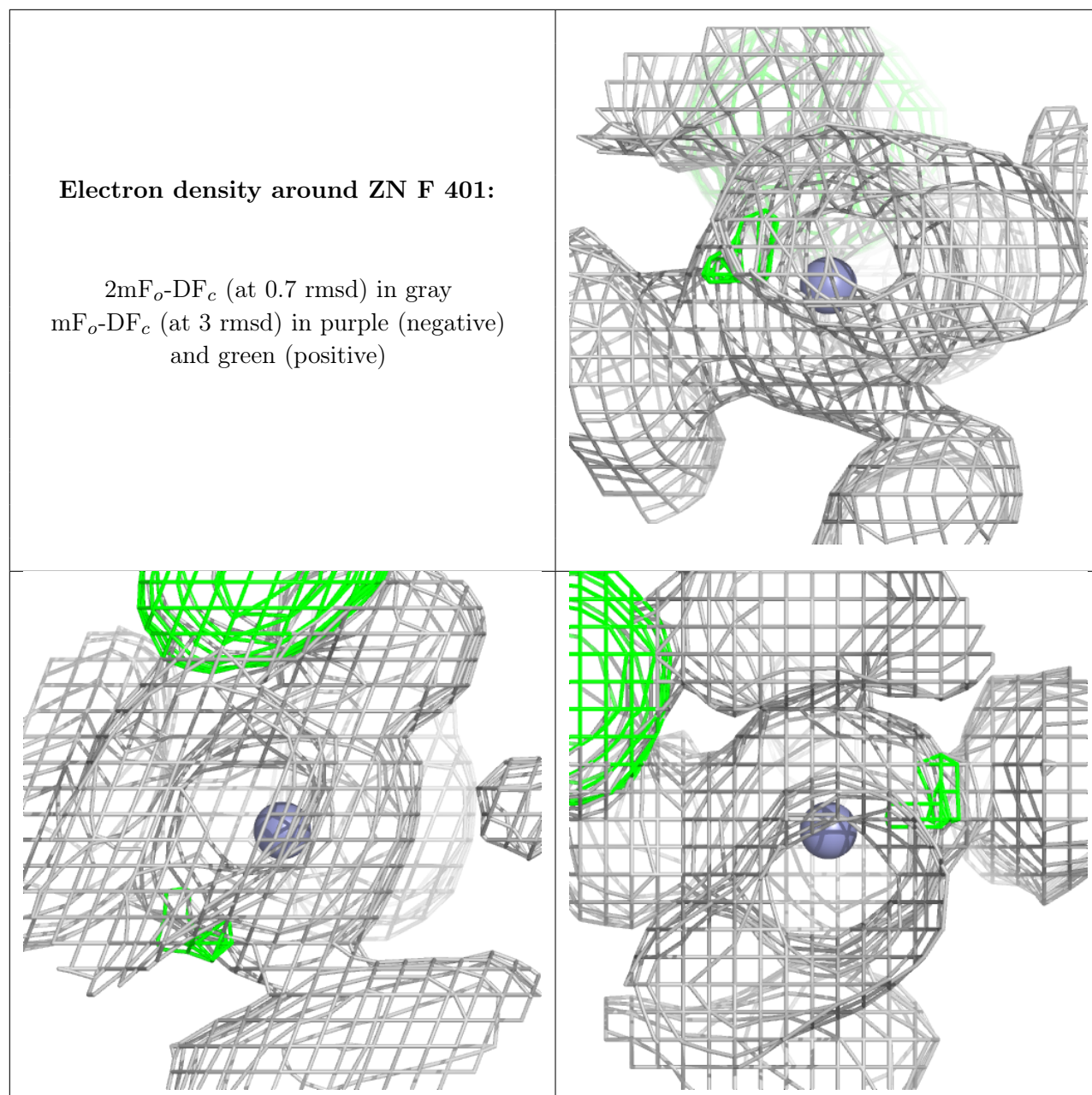
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around ZN C 401:

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





6.5 Other polymers ⓘ

There are no such residues in this entry.