



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 04:00 AM UTC

PDB ID : 7BZ1 / pdb_00007bz1
Title : The mutant variant of PNGM-1. H96 was substituted for alanine to study metal coordination.
Authors : Park, Y.S.; Kang, L.W.; Lee, J.H.
Deposited on : 2020-04-26
Resolution : 2.45 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

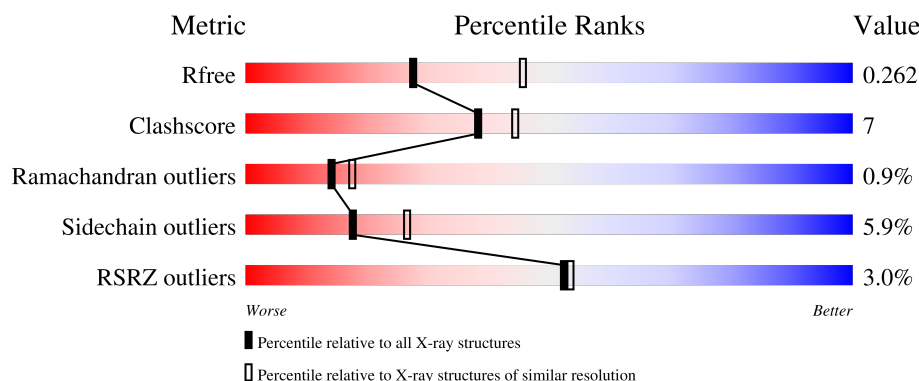
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.45 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	1190 (2.46-2.46)
Clashscore	190562	1229 (2.46-2.46)
Ramachandran outliers	187476	1218 (2.46-2.46)
Sidechain outliers	187428	1218 (2.46-2.46)
RSRZ outliers	180081	1190 (2.46-2.46)

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	365	2% 82% 17% .
1	B	365	2% 81% 17% .
1	C	365	4% 75% 22% .
1	D	365	3% 81% 18% .

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 11849 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	365	Total	C	N	O	S	0	0	0
			2886	1825	496	545	20			
1	B	365	Total	C	N	O	S	0	0	0
			2886	1825	496	545	20			
1	C	365	Total	C	N	O	S	0	0	0
			2882	1823	496	543	20			
1	D	365	Total	C	N	O	S	0	0	0
			2883	1824	496	543	20			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	96	ALA	HIS	engineered mutation	UNP A0A2U8UYM6
B	96	ALA	HIS	engineered mutation	UNP A0A2U8UYM6
C	96	ALA	HIS	engineered mutation	UNP A0A2U8UYM6
D	96	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	0	0
			1	1		
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		

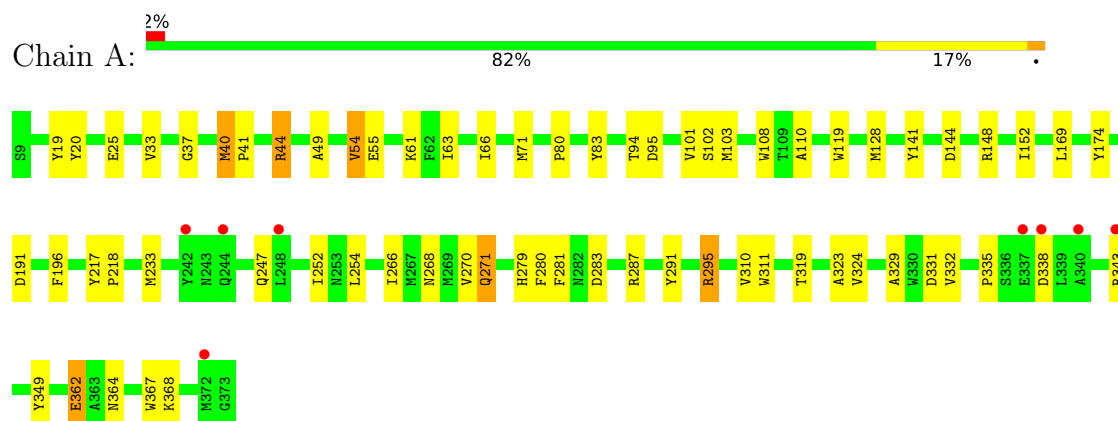
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	80	Total 80	O 80	0	0
3	B	74	Total 74	O 74	0	0
3	C	59	Total 59	O 59	0	0
3	D	95	Total 95	O 95	0	0

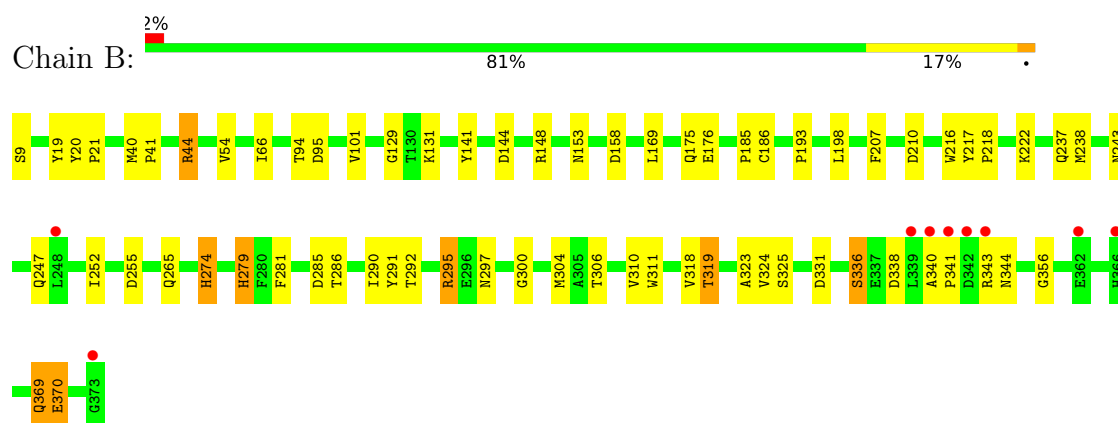
3 Residue-property plots

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.

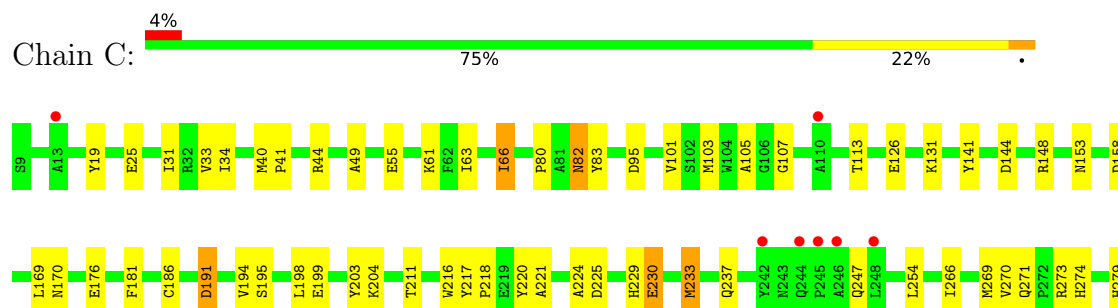
• Molecule 1: Metallo-beta-lactamase PNGM-1



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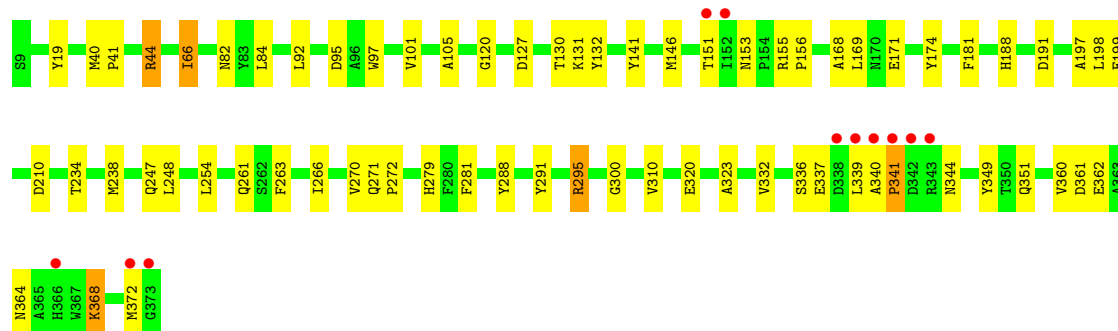
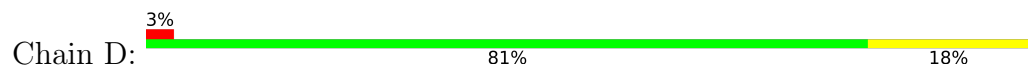


• Molecule 1: Metallo-beta-lactamase PNGM-1





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



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	121.64Å 82.55Å 163.72Å 90.00° 110.77° 90.00°	Depositor
Resolution (Å)	48.08 – 2.45 48.08 – 2.45	Depositor EDS
% Data completeness (in resolution range)	90.0 (48.08-2.45) 90.2 (48.08-2.45)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.63 (at 2.45Å)	Xtriage
Refinement program	REFMAC 5.8.0257	Depositor
R, R_{free}	0.186 , 0.260 0.192 , 0.262	Depositor DCC
R_{free} test set	2485 reflections (4.43%)	wwPDB-VP
Wilson B-factor (Å ²)	25.0	Xtriage
Anisotropy	0.172	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 27.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.008 for h,-k,-h-l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	11849	wwPDB-VP
Average B, all atoms (Å ²)	31.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.91% 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: 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	1.11	1/2971 (0.0%)	1.47	5/4051 (0.1%)
1	B	1.12	1/2971 (0.0%)	1.42	3/4051 (0.1%)
1	C	1.12	1/2967 (0.0%)	1.46	10/4046 (0.2%)
1	D	1.10	1/2968 (0.0%)	1.48	7/4047 (0.2%)
All	All	1.11	4/11877 (0.0%)	1.46	25/16195 (0.2%)

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	271	GLN	N-CA	5.42	1.50	1.46
1	B	274	HIS	CE1-NE2	5.28	1.37	1.32
1	A	271	GLN	N-CA	5.19	1.50	1.46
1	D	84	LEU	C-O	5.09	1.30	1.24

All (25) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	95	ASP	CA-CB-CG	6.65	119.25	112.60
1	D	360	VAL	CA-C-N	6.35	128.70	120.44
1	D	360	VAL	C-N-CA	6.35	128.70	120.44
1	A	66	ILE	CA-C-N	5.48	126.72	120.53
1	A	66	ILE	C-N-CA	5.48	126.72	120.53
1	C	66	ILE	CA-C-N	5.48	126.41	120.44
1	C	66	ILE	C-N-CA	5.48	126.41	120.44
1	D	361	ASP	CA-C-N	5.43	128.56	120.31
1	D	361	ASP	C-N-CA	5.43	128.56	120.31
1	D	263	PHE	CA-C-N	5.33	126.00	120.03
1	D	263	PHE	C-N-CA	5.33	126.00	120.03
1	B	255	ASP	CA-CB-CG	5.29	117.89	112.60
1	A	196	PHE	CA-C-N	5.19	130.09	122.77

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	196	PHE	C-N-CA	5.19	130.09	122.77
1	A	95	ASP	CA-CB-CG	5.16	117.76	112.60
1	B	129	GLY	CA-C-N	5.12	127.14	120.28
1	B	129	GLY	C-N-CA	5.12	127.14	120.28
1	C	107	GLY	CA-C-N	5.08	127.40	120.54
1	C	107	GLY	C-N-CA	5.08	127.40	120.54
1	C	105	ALA	CA-C-N	5.05	125.45	119.94
1	C	105	ALA	C-N-CA	5.05	125.45	119.94
1	C	362	GLU	CA-C-N	5.04	127.54	120.28
1	C	362	GLU	C-N-CA	5.04	127.54	120.28
1	C	191	ASP	CA-C-O	-5.01	115.64	121.15
1	C	34	ILE	CA-C-O	-5.00	115.18	120.53

There are no chirality outliers.

There are no planarity outliers.

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	2886	0	2707	43	0
1	B	2886	0	2707	34	0
1	C	2882	0	2703	49	0
1	D	2883	0	2705	47	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	80	0	0	0	0
3	B	74	0	0	3	0
3	C	59	0	0	1	0
3	D	95	0	0	2	0
All	All	11849	0	10822	154	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 7.

All (154) 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:D:151:THR:HA	3:D:588:HOH:O	1.60	1.01
1:C:221:ALA:O	1:C:270:VAL:HG13	1.82	0.79
1:D:44:ARG:HD3	1:D:281:PHE:CE1	2.21	0.75
1:C:316:ASP:OD1	1:C:316:ASP:N	2.20	0.74
1:A:144:ASP:OD1	1:A:148:ARG:NH1	2.21	0.74
1:B:210:ASP:OD2	3:B:501:HOH:O	2.05	0.74
1:A:55:GLU:OE1	1:A:61:LYS:HE3	1.89	0.71
1:C:295:ARG:NH2	1:C:300:GLY:O	2.23	0.71
1:B:101:VAL:HG21	1:B:141:TYR:CE2	2.28	0.68
1:A:291:TYR:CZ	1:A:295:ARG:HD2	2.30	0.67
1:A:364:ASN:O	1:A:368:LYS:HB2	1.95	0.66
1:C:266:ILE:O	1:C:270:VAL:HG23	1.98	0.64
1:D:340:ALA:O	1:D:341:PRO:O	2.16	0.63
1:C:221:ALA:O	1:C:270:VAL:CG1	2.47	0.63
1:D:295:ARG:NH2	1:D:300:GLY:O	2.31	0.62
1:B:40:MET:HB3	1:B:41:PRO:HD2	1.82	0.61
1:C:364:ASN:O	1:C:368:LYS:HB2	2.00	0.61
1:A:80:PRO:HD2	1:A:83:TYR:HD2	1.67	0.60
1:C:82:ASN:HB2	3:C:515:HOH:O	2.02	0.60
1:A:63:ILE:CD1	1:A:103:MET:HE3	2.32	0.59
1:C:237:GLN:NE2	1:C:286:THR:HA	2.18	0.59
1:C:82:ASN:HD21	1:C:113:THR:HG23	1.68	0.59
1:C:40:MET:HB3	1:C:41:PRO:HD2	1.85	0.58
1:D:364:ASN:O	1:D:368:LYS:HB2	2.04	0.58
1:B:356:GLY:HA3	1:C:216:TRP:CH2	2.39	0.58
1:A:40:MET:HE2	1:A:280:PHE:HA	1.86	0.58
1:C:203:TYR:HA	1:C:225:ASP:OD2	2.03	0.57
1:C:19:TYR:O	1:C:323:ALA:HA	2.05	0.57
1:B:44:ARG:HD3	1:B:281:PHE:CE1	2.40	0.57
1:C:290:ILE:HG22	1:C:304:MET:HE1	1.86	0.56
1:D:198:LEU:C	1:D:198:LEU:HD23	2.31	0.56
1:C:144:ASP:OD1	1:C:148:ARG:NH1	2.39	0.56
1:D:153:ASN:HD22	1:D:339:LEU:HG	1.70	0.56
1:A:101:VAL:HG21	1:A:141:TYR:CE2	2.40	0.56
1:B:295:ARG:NH2	1:B:300:GLY:O	2.36	0.56
1:C:217:TYR:HB3	1:C:218:PRO:HD3	1.88	0.56
1:B:237:GLN:NE2	1:B:286:THR:HA	2.20	0.55
1:B:238:MET:HE2	1:B:252:ILE:HD13	1.89	0.54
1:A:110:ALA:HB1	1:A:331:ASP:OD1	2.07	0.54
1:A:55:GLU:HG2	1:A:61:LYS:HG2	1.90	0.53
1:A:110:ALA:HB1	1:A:331:ASP:CG	2.33	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:33:VAL:HG22	1:A:54:VAL:HG13	1.90	0.53
1:A:349:TYR:HA	1:D:191:ASP:OD1	2.08	0.53
1:B:144:ASP:OD1	1:B:148:ARG:NH1	2.42	0.52
1:C:229:HIS:C	1:C:230:GLU:O	2.53	0.52
1:D:127:ASP:HA	1:D:132:TYR:CG	2.45	0.52
1:D:266:ILE:O	1:D:270:VAL:HG23	2.10	0.52
1:D:291:TYR:CZ	1:D:295:ARG:HD2	2.45	0.52
1:B:279:HIS:NE2	3:B:501:HOH:O	2.27	0.51
1:B:44:ARG:HD2	1:C:329:ALA:O	2.10	0.51
1:A:20:TYR:CD1	1:B:20:TYR:CD1	2.99	0.51
1:C:198:LEU:C	1:C:198:LEU:HD23	2.36	0.51
1:C:229:HIS:O	1:C:230:GLU:C	2.54	0.51
1:A:80:PRO:HD2	1:A:83:TYR:CD2	2.45	0.51
1:C:101:VAL:HG21	1:C:141:TYR:CE2	2.46	0.50
1:A:266:ILE:O	1:A:270:VAL:HG23	2.12	0.50
1:A:287:ARG:HD3	1:C:288:TYR:OH	2.12	0.50
1:B:153:ASN:ND2	1:B:340:ALA:O	2.40	0.50
1:A:63:ILE:HD12	1:A:103:MET:HE3	1.92	0.50
1:C:170:ASN:ND2	1:C:220:TYR:CD1	2.80	0.49
1:D:153:ASN:ND2	1:D:339:LEU:HG	2.27	0.49
1:A:94:THR:HG22	1:D:101:VAL:HG22	1.94	0.49
1:A:283:ASP:O	1:A:287:ARG:HG2	2.13	0.49
1:B:285:ASP:HB3	1:D:320:GLU:OE2	2.13	0.48
1:C:40:MET:HB3	1:C:41:PRO:CD	2.42	0.48
1:A:311:TRP:HA	1:A:319:THR:O	2.14	0.48
1:B:101:VAL:HG21	1:B:141:TYR:CZ	2.48	0.48
1:B:186:CYS:HB2	1:B:193:PRO:HB2	1.95	0.48
1:A:40:MET:HE1	1:A:233:MET:HE2	1.95	0.48
1:B:94:THR:CG2	1:C:101:VAL:HG22	2.44	0.48
1:D:66:ILE:HG23	1:D:66:ILE:O	2.14	0.48
1:D:181:PHE:HA	1:D:197:ALA:O	2.13	0.48
1:D:101:VAL:HG21	1:D:141:TYR:CE2	2.49	0.48
1:C:186:CYS:HB3	1:C:211:THR:HB	1.96	0.47
1:D:234:THR:O	1:D:238:MET:HG2	2.13	0.47
1:D:368:LYS:O	1:D:372:MET:SD	2.72	0.47
1:D:44:ARG:HD3	1:D:281:PHE:CZ	2.50	0.47
1:D:40:MET:HB3	1:D:41:PRO:HD2	1.96	0.47
1:C:153:ASN:HB2	1:C:340:ALA:O	2.15	0.47
1:A:191:ASP:OD1	1:D:349:TYR:HA	2.15	0.47
1:B:290:ILE:HG22	1:B:304:MET:HE1	1.96	0.47
1:D:153:ASN:HB2	1:D:341:PRO:HA	1.97	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:217:TYR:HB3	1:B:218:PRO:HD3	1.97	0.46
1:C:198:LEU:HD23	1:C:199:GLU:N	2.30	0.46
1:C:204:LYS:HB3	1:C:224:ALA:HA	1.96	0.46
1:B:185:PRO:HB2	1:B:216:TRP:CZ3	2.51	0.46
1:B:54:VAL:HG21	1:B:207:PHE:CZ	2.50	0.46
1:C:49:ALA:HB1	1:C:95:ASP:HB2	1.97	0.46
1:A:128:MET:HE1	1:D:146:MET:HB2	1.98	0.45
1:B:274:HIS:CB	1:B:318:VAL:HG11	2.46	0.45
1:C:274:HIS:CB	1:C:318:VAL:HG11	2.47	0.45
1:C:283:ASP:O	1:C:285:ASP:N	2.50	0.45
1:B:198:LEU:C	1:B:198:LEU:HD23	2.42	0.45
1:A:94:THR:CG2	1:D:101:VAL:HG22	2.47	0.45
1:C:283:ASP:C	1:C:285:ASP:H	2.25	0.44
1:D:247:GLN:NE2	1:D:248:LEU:CD1	2.81	0.44
1:A:108:TRP:CE3	1:A:148:ARG:HD2	2.52	0.44
1:D:155:ARG:HB3	1:D:156:PRO:HD3	1.98	0.44
1:A:119:TRP:CZ2	1:A:174:TYR:HD2	2.35	0.44
1:A:362:GLU:H	1:A:362:GLU:CD	2.26	0.44
1:B:94:THR:HG22	1:C:101:VAL:HG22	2.00	0.44
1:B:265:GLN:OE1	1:C:360:VAL:HA	2.18	0.44
1:C:44:ARG:HD3	1:C:281:PHE:CE1	2.53	0.44
1:A:268:ASN:O	1:A:271:GLN:NE2	2.51	0.43
1:C:225:ASP:HA	1:C:273:ARG:HD3	2.01	0.43
1:A:40:MET:HB3	1:A:41:PRO:HD2	1.99	0.43
1:C:63:ILE:CD1	1:C:103:MET:HE3	2.48	0.43
1:B:291:TYR:CZ	1:D:288:TYR:HB3	2.54	0.43
1:C:218:PRO:HB3	1:C:269:MET:HE2	2.01	0.43
1:A:37:GLY:HA3	1:A:49:ALA:O	2.19	0.43
1:D:168:ALA:HB1	1:D:171:GLU:CG	2.49	0.43
1:D:279:HIS:CE1	3:D:501:HOH:O	2.71	0.43
1:A:367:TRP:CD1	1:D:261:GLN:CD	2.97	0.43
1:A:83:TYR:OH	1:A:335:PRO:HG3	2.19	0.42
1:A:324:VAL:HG12	1:B:324:VAL:HG12	2.01	0.42
1:D:40:MET:HB3	1:D:41:PRO:CD	2.48	0.42
1:C:186:CYS:SG	1:C:194:VAL:C	3.03	0.42
1:D:188:HIS:CD2	1:D:210:ASP:HB2	2.55	0.42
1:B:369:GLN:HG2	1:B:370:GLU:N	2.35	0.42
1:C:33:VAL:O	1:C:310:VAL:HA	2.19	0.42
1:B:19:TYR:O	1:B:323:ALA:HA	2.20	0.42
1:C:291:TYR:CZ	1:C:295:ARG:HD2	2.55	0.42
1:C:291:TYR:CE1	1:C:302:LEU:HD23	2.55	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:82:ASN:HD22	1:D:336:SER:HB3	1.85	0.42
1:D:153:ASN:OD1	1:D:153:ASN:C	2.63	0.42
1:A:94:THR:HG21	1:D:105:ALA:HB2	2.02	0.42
1:B:311:TRP:HA	1:B:319:THR:O	2.19	0.42
1:D:40:MET:CG	1:D:281:PHE:CE1	3.03	0.42
1:A:19:TYR:O	1:A:323:ALA:HA	2.20	0.42
3:B:547:HOH:O	1:C:141:TYR:HA	2.19	0.42
1:D:19:TYR:O	1:D:323:ALA:HA	2.20	0.42
1:D:247:GLN:NE2	1:D:248:LEU:HD13	2.35	0.42
1:A:44:ARG:NH2	1:A:283:ASP:OD1	2.32	0.41
1:A:329:ALA:O	1:D:44:ARG:HD2	2.20	0.41
1:C:55:GLU:HG2	1:C:61:LYS:HE3	2.01	0.41
1:D:198:LEU:HD23	1:D:199:GLU:N	2.35	0.41
1:B:54:VAL:CG2	1:B:207:PHE:HZ	2.33	0.41
1:D:198:LEU:C	1:D:198:LEU:CD2	2.93	0.41
1:A:217:TYR:HB3	1:A:218:PRO:HD3	2.03	0.41
1:C:31:ILE:HB	1:C:313:ILE:HB	2.02	0.41
1:C:80:PRO:HD2	1:C:83:TYR:HD2	1.86	0.41
1:D:174:TYR:HB3	1:D:181:PHE:HB2	2.01	0.41
1:A:61:LYS:NZ	1:B:9:SER:O	2.43	0.41
1:A:148:ARG:HG3	1:A:152:ILE:HD12	2.02	0.41
1:D:92:LEU:HA	1:D:97:TRP:CD1	2.56	0.41
1:A:44:ARG:HD3	1:A:281:PHE:CE1	2.56	0.41
1:B:54:VAL:HG21	1:B:207:PHE:HZ	1.86	0.41
1:D:271:GLN:N	1:D:272:PRO:CD	2.84	0.41
1:A:71:MET:HG2	1:A:102:SER:OG	2.21	0.40
1:D:168:ALA:HB1	1:D:171:GLU:HG3	2.03	0.40
1:C:181:PHE:CZ	1:C:198:LEU:HD12	2.57	0.40
1:D:120:GLY:HA3	1:D:130:THR:HG21	2.01	0.40
1:B:21:PRO:HG2	1:B:325:SER:HB3	2.04	0.40
1:C:233:MET:HE2	1:C:233:MET:HB2	1.83	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	363/365 (100%)	337 (93%)	25 (7%)	1 (0%)	36	45
1	B	363/365 (100%)	340 (94%)	18 (5%)	5 (1%)	9	8
1	C	363/365 (100%)	330 (91%)	28 (8%)	5 (1%)	9	8
1	D	363/365 (100%)	335 (92%)	26 (7%)	2 (1%)	21	27
All	All	1452/1460 (100%)	1342 (92%)	97 (7%)	13 (1%)	14	17

All (13) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	341	PRO
1	D	341	PRO
1	B	341	PRO
1	C	284	ASP
1	B	279	HIS
1	C	230	GLU
1	A	279	HIS
1	C	336	SER
1	B	243	ASN
1	B	336	SER
1	C	66	ILE
1	D	66	ILE
1	B	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	297/297 (100%)	283 (95%)	14 (5%)	23	35
1	B	297/297 (100%)	275 (93%)	22 (7%)	13	16
1	C	296/297 (100%)	274 (93%)	22 (7%)	13	16
1	D	296/297 (100%)	284 (96%)	12 (4%)	27	40

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
All	All	1186/1188 (100%)	1116 (94%)	70 (6%)	18	26

All (70) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	25	GLU
1	A	40	MET
1	A	44	ARG
1	A	54	VAL
1	A	169	LEU
1	A	247	GLN
1	A	252	ILE
1	A	254	LEU
1	A	295	ARG
1	A	310	VAL
1	A	332	VAL
1	A	338	ASP
1	A	343	ARG
1	A	362	GLU
1	B	44	ARG
1	B	95	ASP
1	B	131	LYS
1	B	158	ASP
1	B	169	LEU
1	B	175	GLN
1	B	176	GLU
1	B	222	LYS
1	B	247	GLN
1	B	292	THR
1	B	295	ARG
1	B	297	ASN
1	B	306	THR
1	B	310	VAL
1	B	319	THR
1	B	331	ASP
1	B	336	SER
1	B	338	ASP
1	B	343	ARG
1	B	344	ASN
1	B	369	GLN
1	B	370	GLU
1	C	25	GLU

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Mol	Chain	Res	Type
1	C	82	ASN
1	C	126	GLU
1	C	131	LYS
1	C	158	ASP
1	C	169	LEU
1	C	176	GLU
1	C	191	ASP
1	C	195	SER
1	C	233	MET
1	C	247	GLN
1	C	254	LEU
1	C	279	HIS
1	C	295	ARG
1	C	310	VAL
1	C	316	ASP
1	C	332	VAL
1	C	336	SER
1	C	337	GLU
1	C	338	ASP
1	C	367	TRP
1	C	368	LYS
1	D	44	ARG
1	D	131	LYS
1	D	169	LEU
1	D	254	LEU
1	D	295	ARG
1	D	310	VAL
1	D	332	VAL
1	D	337	GLU
1	D	344	ASN
1	D	351	GLN
1	D	362	GLU
1	D	368	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (16) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	82	ASN
1	A	175	GLN
1	A	237	GLN
1	A	257	HIS
1	A	359	ASN

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Mol	Chain	Res	Type
1	A	369	GLN
1	B	82	ASN
1	B	170	ASN
1	B	175	GLN
1	B	344	ASN
1	C	75	GLN
1	C	82	ASN
1	C	237	GLN
1	D	82	ASN
1	D	175	GLN
1	D	369	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 4 ligands modelled in this entry, 4 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	365/365 (100%)	-0.07	8 (2%) 62 64	10, 21, 67, 109	0
1	B	365/365 (100%)	0.03	9 (2%) 58 61	14, 25, 70, 114	0
1	C	365/365 (100%)	0.24	16 (4%) 39 38	15, 30, 69, 110	0
1	D	365/365 (100%)	0.08	11 (3%) 52 53	13, 26, 65, 137	0
All	All	1460/1460 (100%)	0.07	44 (3%) 52 53	10, 26, 68, 137	0

All (44) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	342	ASP	6.2
1	D	339	LEU	5.6
1	D	152	ILE	4.6
1	D	340	ALA	4.4
1	C	340	ALA	3.9
1	C	339	LEU	3.7
1	D	341	PRO	3.5
1	B	339	LEU	3.4
1	B	340	ALA	3.2
1	B	343	ARG	3.1
1	D	373	GLY	3.1
1	B	341	PRO	3.0
1	C	248	LEU	3.0
1	D	151	THR	2.9
1	D	343	ARG	2.8
1	A	337	GLU	2.8
1	A	338	ASP	2.7
1	C	373	GLY	2.7
1	C	341	PRO	2.6
1	C	343	ARG	2.6
1	A	242	TYR	2.5

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Mol	Chain	Res	Type	RSRZ
1	B	373	GLY	2.5
1	D	366	HIS	2.5
1	C	332	VAL	2.5
1	C	338	ASP	2.4
1	C	110	ALA	2.4
1	B	248	LEU	2.4
1	C	244	GLN	2.3
1	B	362	GLU	2.3
1	C	242	TYR	2.3
1	A	372	MET	2.3
1	C	13	ALA	2.2
1	B	366	HIS	2.2
1	A	248	LEU	2.2
1	A	340	ALA	2.2
1	B	342	ASP	2.2
1	A	343	ARG	2.2
1	D	372	MET	2.1
1	C	246	ALA	2.1
1	C	337	GLU	2.1
1	D	338	ASP	2.1
1	C	245	PRO	2.1
1	A	244	GLN	2.0
1	C	365	ALA	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

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

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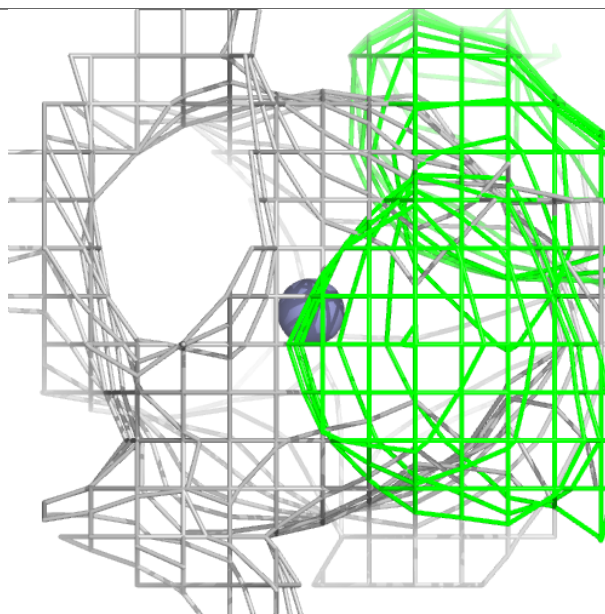
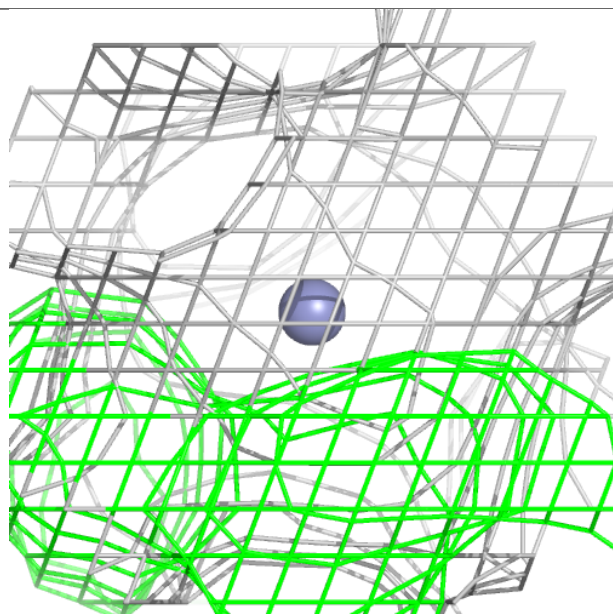
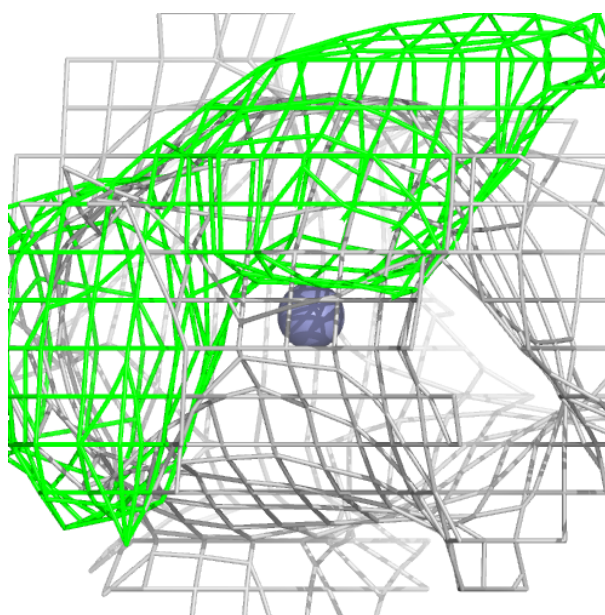
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	ZN	A	401	1/1	0.99	0.03	22,22,22,22	0
2	ZN	B	401	1/1	0.99	0.05	26,26,26,26	0
2	ZN	C	401	1/1	0.99	0.04	30,30,30,30	0
2	ZN	D	401	1/1	0.99	0.04	24,24,24,24	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.

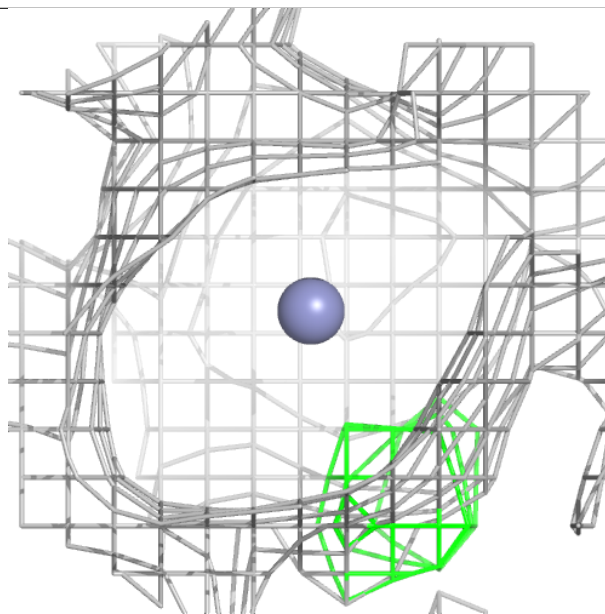
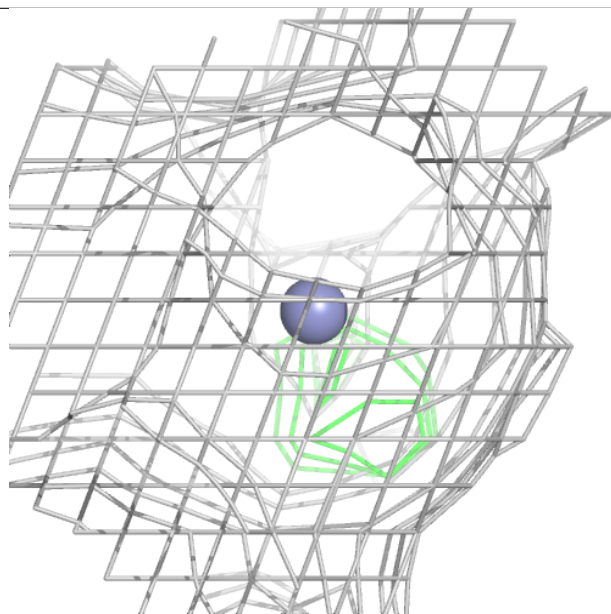
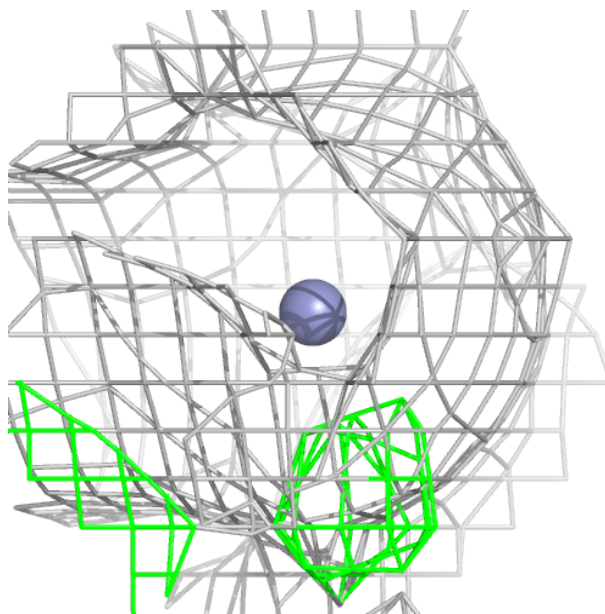
Electron density around ZN A 401:

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



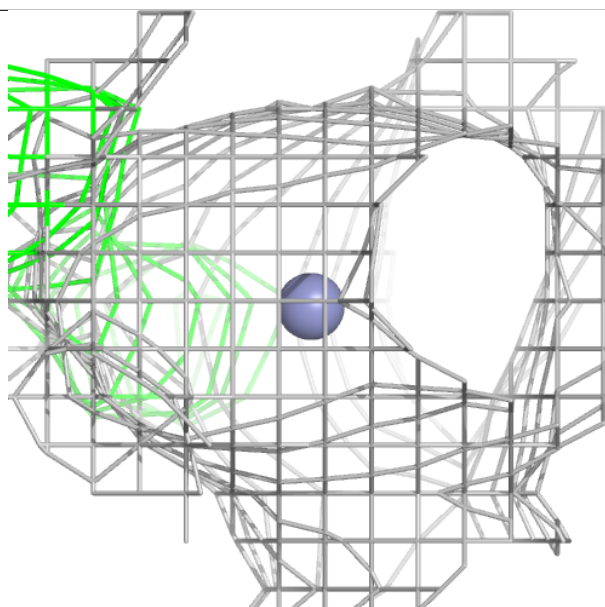
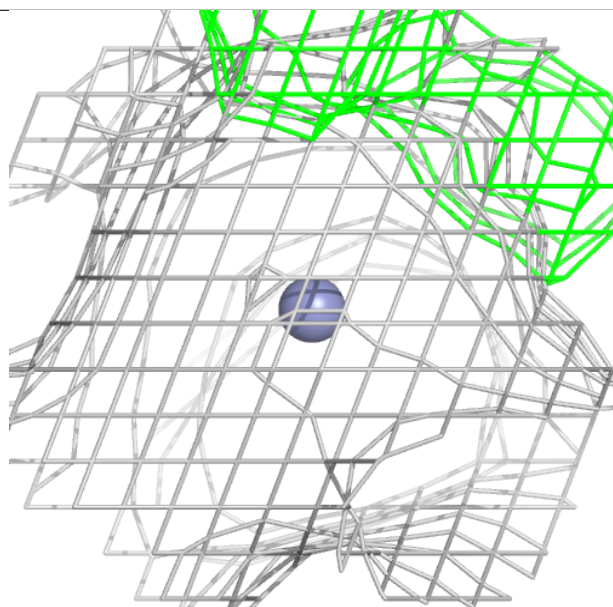
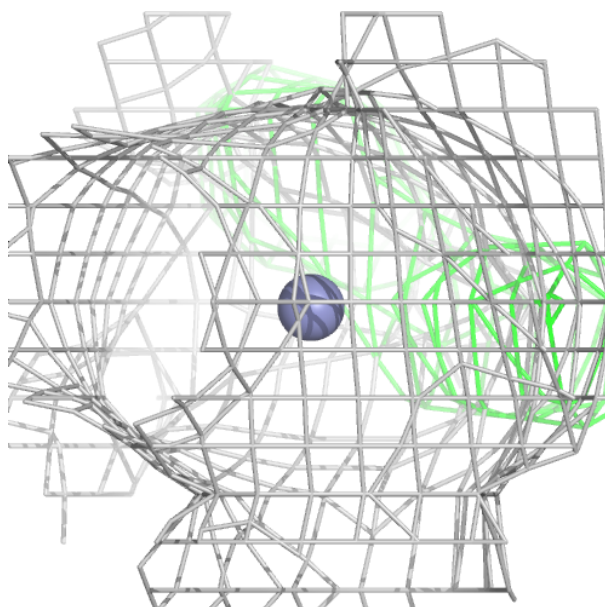
Electron density around ZN B 401:

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



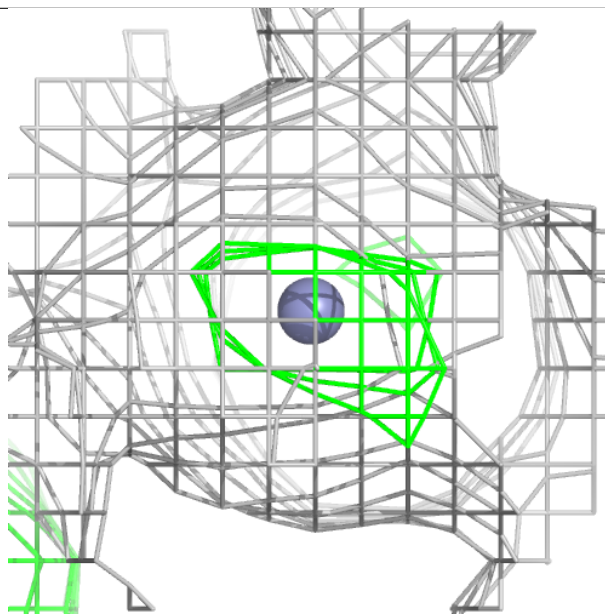
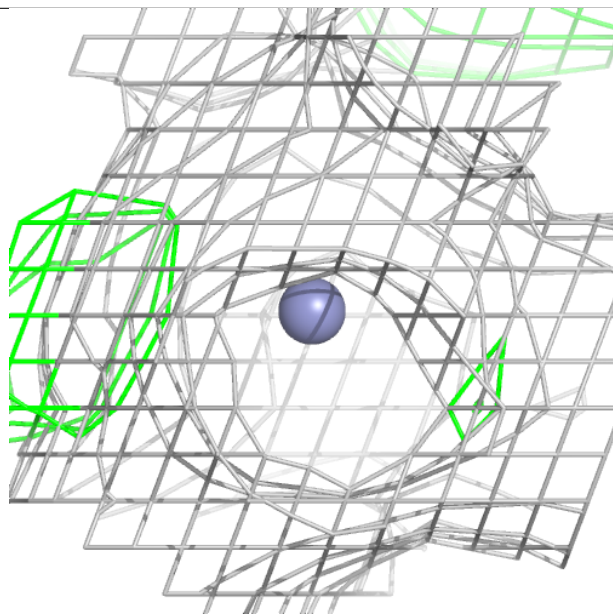
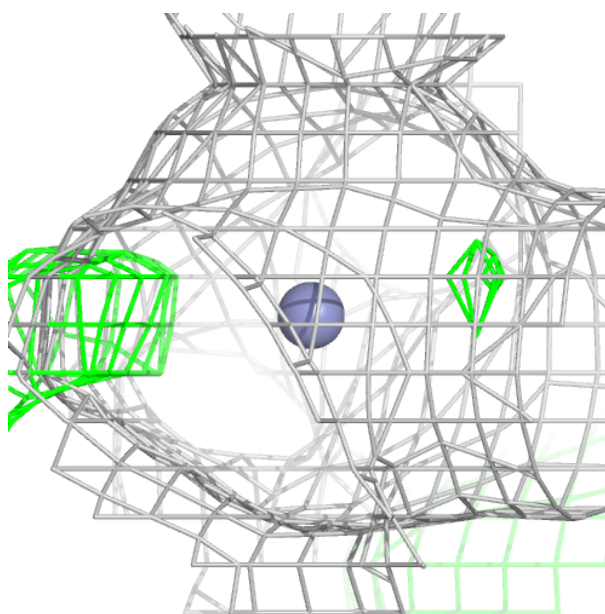
Electron density around ZN C 401:

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



Electron density around ZN D 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.