



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 11:00 PM UTC

PDB ID : 9GN1 / pdb_00009gn1
Title : Crystal structure of inactive Deacetylase (HdaH) H144A from *Klebsiella pneumoniae* subsp. *ozaenae*
Authors : Qin, Q.; Graf, L.G.; Schulze, S.; Palm, G.J.; Lammers, M.
Deposited on : 2024-08-30
Resolution : 2.35 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

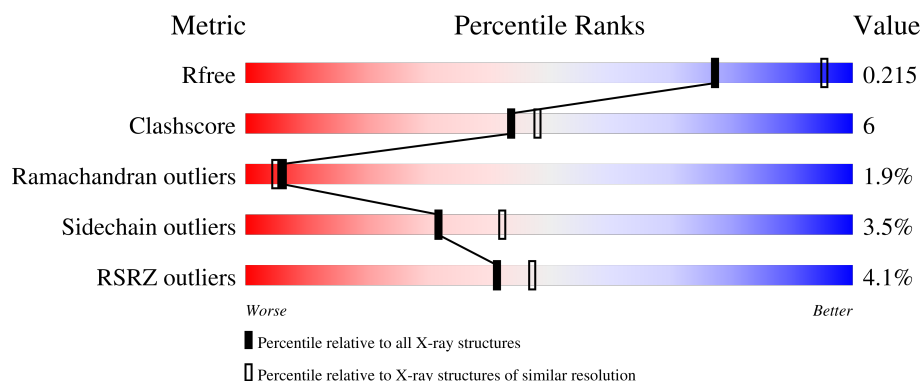
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.35 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	1596 (2.36-2.36)
Clashscore	190562	1663 (2.36-2.36)
Ramachandran outliers	187476	1646 (2.36-2.36)
Sidechain outliers	187428	1646 (2.36-2.36)
RSRZ outliers	180081	1598 (2.36-2.36)

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	382	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
4	IMD	A	405	-	-	X	-

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Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
5	ACT	A	406	-	X	X	-

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 5735 atoms, of which 2768 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 Deacetylase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	370	Total	C	H	N	O	S	69	0	0
			5577	1775	2760	499	525	18			

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-10	MET	-	initiating methionine	UNP A0A377Z5F6
A	-9	ALA	-	expression tag	UNP A0A377Z5F6
A	-8	HIS	-	expression tag	UNP A0A377Z5F6
A	-7	HIS	-	expression tag	UNP A0A377Z5F6
A	-6	HIS	-	expression tag	UNP A0A377Z5F6
A	-5	HIS	-	expression tag	UNP A0A377Z5F6
A	-4	HIS	-	expression tag	UNP A0A377Z5F6
A	-3	HIS	-	expression tag	UNP A0A377Z5F6
A	-2	VAL	-	expression tag	UNP A0A377Z5F6
A	-1	GLY	-	expression tag	UNP A0A377Z5F6
A	0	THR	-	expression tag	UNP A0A377Z5F6
A	144	ALA	HIS	engineered mutation	UNP A0A377Z5F6

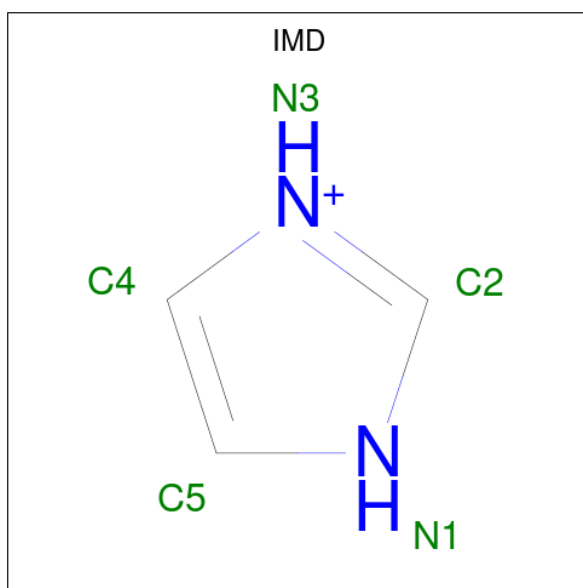
- Molecule 2 is POTASSIUM ION (CCD ID: K) (formula: K).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	2	Total	K	0	0
			2	2		

- Molecule 3 is ZINC ION (CCD ID: ZN) (formula: Zn) (labeled as "Ligand of Interest" by depositor).

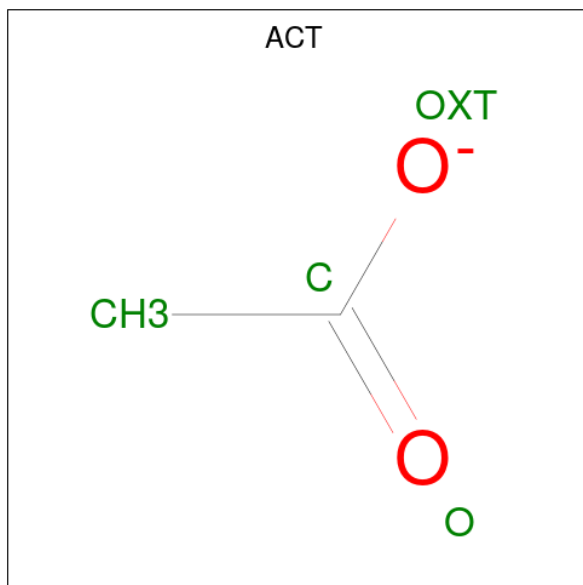
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	2	Total	Zn	0	0
			2	2		

- Molecule 4 is IMIDAZOLE (CCD ID: IMD) (formula: $C_3H_5N_2$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	1	Total	C	H	N	0	0
			10	3	5	2		

- Molecule 5 is ACETATE ION (CCD ID: ACT) (formula: $C_2H_3O_2$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	A	1	Total	C	H	O	3	0
			7	2	3	2		

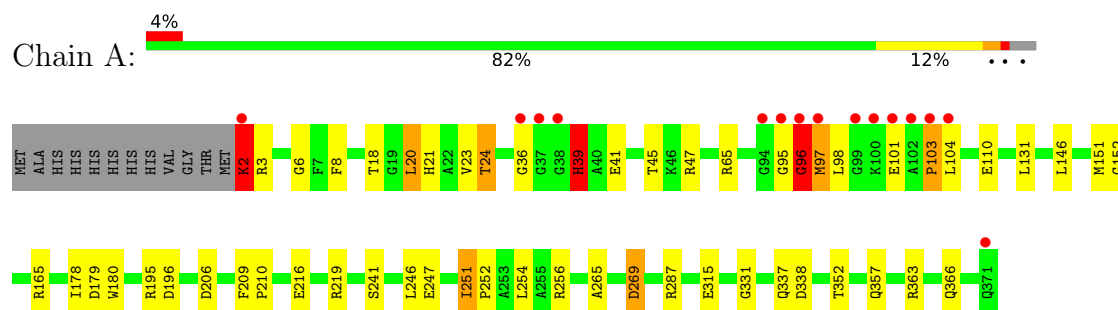
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	137	Total 137	O 137	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: Deacetylase



4 Data and refinement statistics

Property	Value	Source
Space group	I 2 3	Depositor
Cell constants a, b, c, α , β , γ	146.12Å 146.12Å 146.12Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	46.25 – 2.35 46.25 – 2.35	Depositor EDS
% Data completeness (in resolution range)	100.0 (46.25-2.35) 100.0 (46.25-2.35)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.51 (at 2.34Å)	Xtriage
Refinement program	REFMAC 5.8.0425, REFMAC 5.8.0425	Depositor
R, R_{free}	0.153 , 0.204 0.158 , 0.215	Depositor DCC
R_{free} test set	1152 reflections (5.29%)	wwPDB-VP
Wilson B-factor (Å ²)	42.4	Xtriage
Anisotropy	0.000	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.43 , 42.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.032 for -l,-k,-h	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	5735	wwPDB-VP
Average B, all atoms (Å ²)	47.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.93% 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: K, ZN, IMD, ACT

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	0.92	0/2882	1.45	19/3907 (0.5%)

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	A	0	4

There are no bond length outliers.

All (19) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	39	HIS	N-CA-CB	6.98	122.29	110.49
1	A	352	THR	CA-CB-OG1	-6.72	99.52	109.60
1	A	45	THR	CA-CB-OG1	6.67	119.61	109.60
1	A	363	ARG	CB-CA-C	6.29	122.75	110.67
1	A	315	GLU	CB-CG-CD	6.14	123.04	112.60
1	A	39	HIS	CA-CB-CG	-6.11	107.69	113.80
1	A	269	ASP	CA-CB-CG	6.07	118.67	112.60
1	A	24	THR	N-CA-CB	-6.00	101.83	110.65
1	A	196	ASP	CA-CB-CG	5.99	118.59	112.60
1	A	366	GLN	N-CA-CB	5.82	118.44	110.01
1	A	287	ARG	NE-CZ-NH1	-5.51	115.99	121.50
1	A	338	ASP	CA-CB-CG	5.47	118.07	112.60
1	A	216	GLU	N-CA-CB	5.43	118.72	110.14
1	A	206	ASP	CA-CB-CG	5.42	118.02	112.60
1	A	47	ARG	NE-CZ-NH2	5.39	124.05	119.20
1	A	96	GLY	CA-C-N	5.34	131.74	121.54

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	96	GLY	C-N-CA	5.34	131.74	121.54
1	A	110	GLU	CG-CD-OE2	-5.29	106.24	118.40
1	A	195	ARG	NE-CZ-NH2	5.03	123.72	119.20

There are no chirality outliers.

All (4) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	165	ARG	Sidechain
1	A	2	LYS	Peptide
1	A	219	ARG	Sidechain
1	A	256	ARG	Sidechain

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	2817	2760	2749	26	0
2	A	2	0	0	0	0
3	A	2	0	0	0	0
4	A	5	5	5	5	0
5	A	4	3	3	8	0
6	A	137	0	0	4	0
All	All	2967	2768	2757	34	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 6.

All (34) 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:152:GLY:O	6:A:501:HOH:O	1.98	0.81
4:A:405:IMD:HN3	5:A:406:ACT:H2	1.57	0.70
5:A:406:ACT:C	6:A:501:HOH:O	2.49	0.61
4:A:405:IMD:N3	5:A:406:ACT:H2	2.16	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:21:HIS:CG	1:A:39:HIS:HD2	2.20	0.59
5:A:406:ACT:CH3	6:A:501:HOH:O	2.52	0.57
1:A:21:HIS:CG	1:A:39:HIS:CD2	2.93	0.56
1:A:96:GLY:O	1:A:97:MET:HG2	2.04	0.56
1:A:2:LYS:HE3	1:A:331:GLY:HA3	1.87	0.56
1:A:8:PHE:HA	1:A:65:ARG:O	2.09	0.53
1:A:21:HIS:CD2	1:A:39:HIS:HD2	2.28	0.52
4:A:405:IMD:H2	5:A:406:ACT:CH3	2.40	0.52
1:A:21:HIS:HB2	1:A:39:HIS:HB2	1.91	0.51
1:A:146:LEU:HD11	1:A:151:MET:HE1	1.94	0.50
4:A:405:IMD:C2	5:A:406:ACT:CH3	2.90	0.49
1:A:251:ILE:HD12	1:A:254:LEU:HD12	1.95	0.48
1:A:146:LEU:CD1	1:A:151:MET:HE1	2.43	0.48
1:A:2:LYS:HE2	1:A:3:ARG:H	1.79	0.48
1:A:241:SER:HB3	1:A:357:GLN:HG2	1.96	0.48
1:A:18:THR:HG22	1:A:103:PRO:O	2.15	0.47
1:A:20:LEU:HD23	1:A:20:LEU:HA	1.80	0.45
1:A:178:ILE:HD11	1:A:254:LEU:HD11	1.97	0.45
5:A:406:ACT:H3	6:A:501:HOH:O	2.17	0.45
1:A:179:ASP:HA	1:A:265:ALA:HB3	1.99	0.45
1:A:209:PHE:HA	1:A:210:PRO:C	2.42	0.44
1:A:2:LYS:HB3	1:A:2:LYS:NZ	2.33	0.43
1:A:6:GLY:HA3	1:A:131:LEU:HD13	2.01	0.42
1:A:180:TRP:C	1:A:180:TRP:CD1	2.97	0.42
1:A:39:HIS:ND1	1:A:39:HIS:C	2.78	0.41
1:A:98:LEU:O	1:A:98:LEU:HG	2.20	0.41
1:A:251:ILE:HB	1:A:252:PRO:HD3	2.00	0.41
4:A:405:IMD:C2	5:A:406:ACT:H2	2.50	0.40
1:A:95:GLY:O	1:A:96:GLY:C	2.64	0.40
1:A:21:HIS:CB	1:A:39:HIS:CD2	3.04	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	368/382 (96%)	345 (94%)	16 (4%)	7 (2%)	6	5

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	39	HIS
1	A	23	VAL
1	A	96	GLY
1	A	97	MET
1	A	36	GLY
1	A	101	GLU
1	A	103	PRO

5.3.2 Protein sidechains [i](#)

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	288/298 (97%)	278 (96%)	10 (4%)	32	42

All (10) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	2	LYS
1	A	20	LEU
1	A	24	THR
1	A	41	GLU
1	A	104	LEU
1	A	246	LEU
1	A	247	GLU
1	A	251	ILE
1	A	269	ASP
1	A	337	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such

sidechains are listed below:

Mol	Chain	Res	Type
1	A	21	HIS

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 6 ligands modelled in this entry, 4 are monoatomic - leaving 2 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	$\# Z > 2$
4	IMD	A	405	-	5,5,5	0.73	0	5,5,5	0.75	0
5	ACT	A	406	3	3,3,3	2.10	2 (66%)	3,3,3	2.03	1 (33%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	IMD	A	405	-	-	-	0/1/1/1

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	A	406	ACT	CH3-C	2.49	1.58	1.49
5	A	406	ACT	O-C	2.41	1.32	1.22

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	A	406	ACT	O-C-CH3	-2.80	111.06	122.53

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 8 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	405	IMD	5	0
5	A	406	ACT	8	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data [i](#)

6.1 Protein, DNA and RNA chains [i](#)

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2		OWAB(Å ²)	Q<0.9
1	A	370/382 (96%)	-0.49	15 (4%)	41 47	31, 43, 72, 140	0

All (15) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	96	GLY	8.8
1	A	100	LYS	5.1
1	A	101	GLU	4.4
1	A	38	GLY	3.6
1	A	94	GLY	3.0
1	A	99	GLY	3.0
1	A	37	GLY	2.9
1	A	36	GLY	2.9
1	A	2	LYS	2.7
1	A	104	LEU	2.5
1	A	102	ALA	2.5
1	A	371	GLN	2.5
1	A	95	GLY	2.3
1	A	103	PRO	2.2
1	A	97	MET	2.2

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

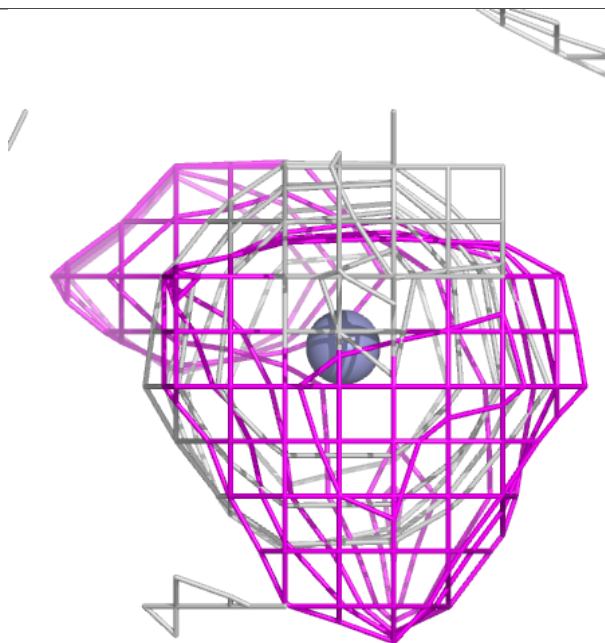
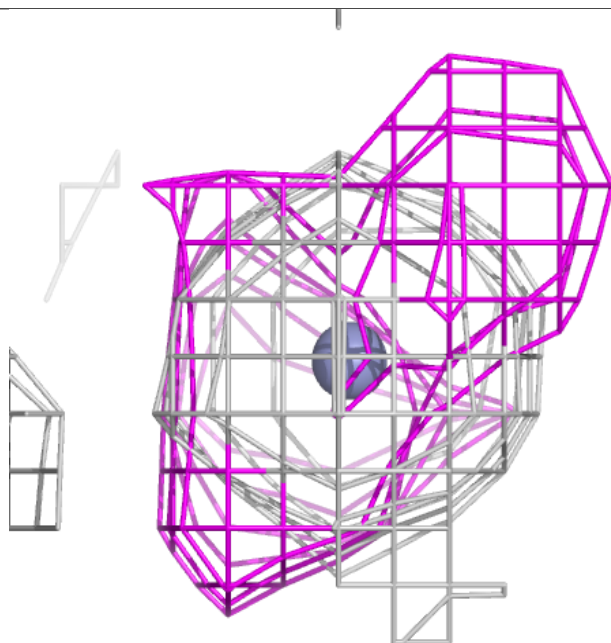
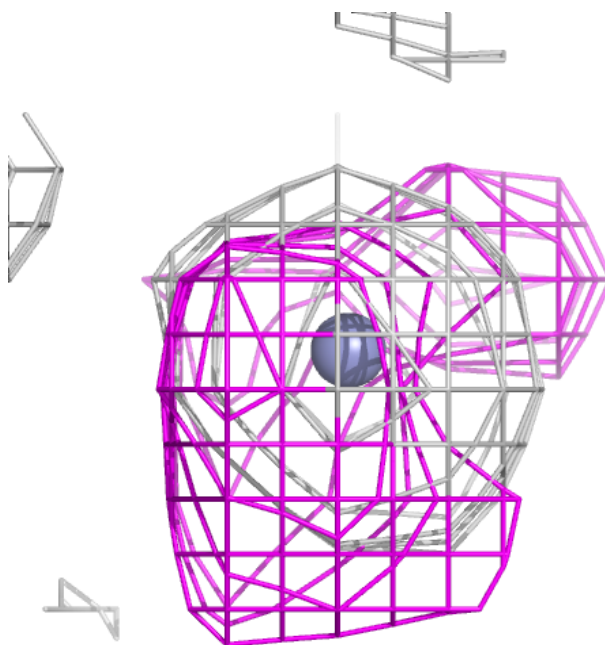
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists 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
4	IMD	A	405	5/5	0.82	0.30	50,62,82,92	0
5	ACT	A	406	4/4	0.95	0.13	28,34,64,72	3
2	K	A	402	1/1	0.99	0.03	55,55,55,55	0
3	ZN	A	404	1/1	0.99	0.21	78,78,78,78	0
3	ZN	A	403	1/1	1.00	0.01	41,41,41,41	0
2	K	A	401	1/1	1.00	0.07	43,43,43,43	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.

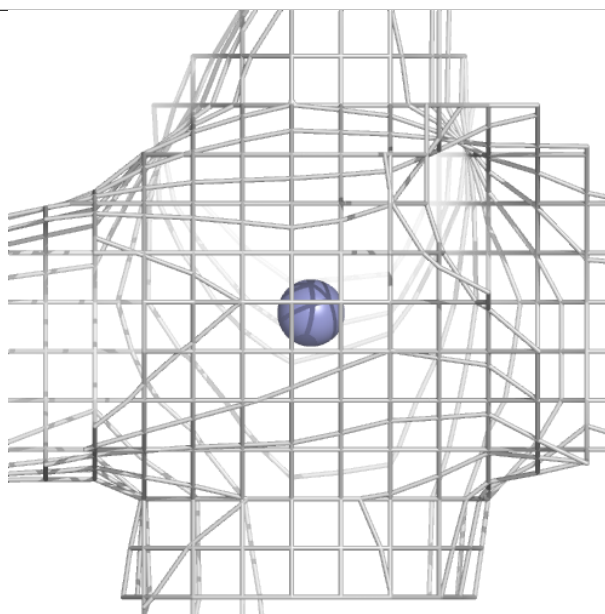
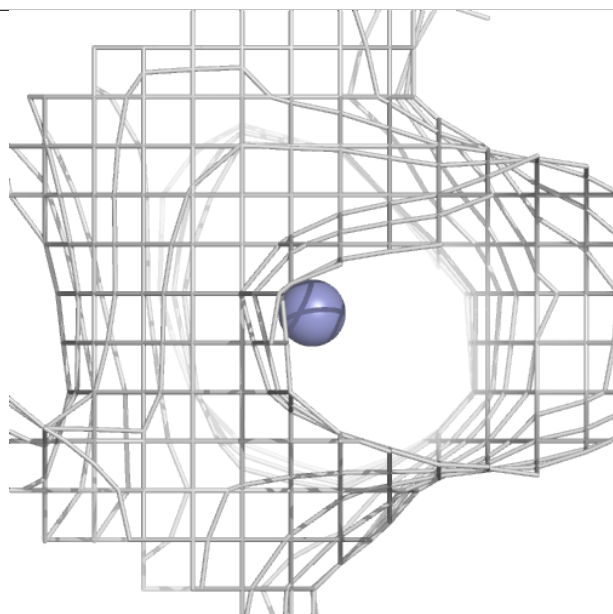
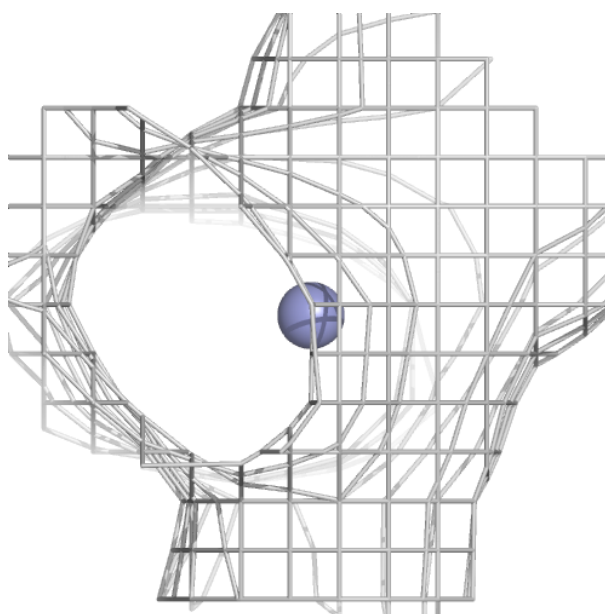
Electron density around ZN A 404:

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



Electron density around ZN A 403:

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



6.5 Other polymers [i](#)

There are no such residues in this entry.