



Full wwPDB X-ray Structure Validation Report ⓘ

Jun 22, 2024 – 11:27 AM EDT

PDB ID : 6K02
Title : Crystal structure of ceNAP1 core
Authors : Liu, Y.R.
Deposited on : 2019-05-05
Resolution : 2.16 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 : 4.02b-467
Xtriage (Phenix) : 1.20.1
EDS : 2.37.1
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.37.1

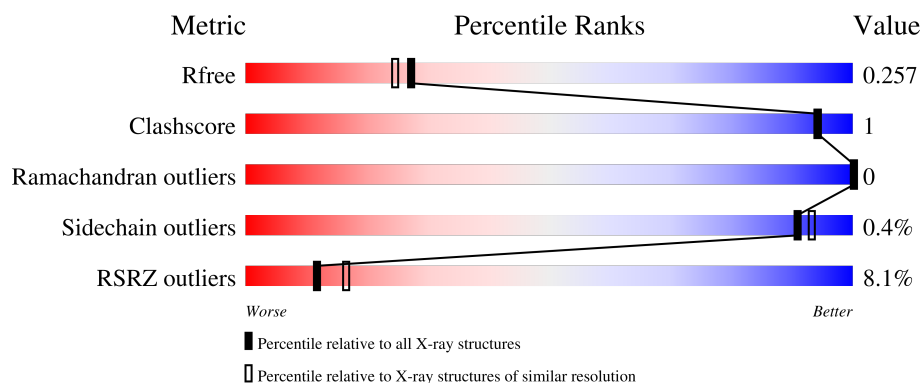
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.16 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}	130704	1479 (2.16-2.16)
Clashscore	141614	1585 (2.16-2.16)
Ramachandran outliers	138981	1560 (2.16-2.16)
Sidechain outliers	138945	1559 (2.16-2.16)
RSRZ outliers	127900	1456 (2.16-2.16)

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	308	<div> <div>7%</div> <div>81%</div> <div>5%</div> <div>15%</div> </div>
1	B	308	<div> <div>7%</div> <div>85%</div> <div>•</div> <div>13%</div> </div>

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 4299 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 Nucleosome Assembly Protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	263	Total	C	N	O	S	0	0	0
			2087	1348	334	401	4			
1	B	269	Total	C	N	O	S	0	0	0
			2149	1382	343	420	4			

There are 42 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-11	MET	-	initiating methionine	UNP Q19007
A	-10	GLY	-	expression tag	UNP Q19007
A	-9	SER	-	expression tag	UNP Q19007
A	-8	SER	-	expression tag	UNP Q19007
A	-7	HIS	-	expression tag	UNP Q19007
A	-6	HIS	-	expression tag	UNP Q19007
A	-5	HIS	-	expression tag	UNP Q19007
A	-4	HIS	-	expression tag	UNP Q19007
A	-3	HIS	-	expression tag	UNP Q19007
A	-2	HIS	-	expression tag	UNP Q19007
A	-1	SER	-	expression tag	UNP Q19007
A	0	SER	-	expression tag	UNP Q19007
A	1	GLY	-	expression tag	UNP Q19007
A	2	GLU	-	expression tag	UNP Q19007
A	3	ASN	-	expression tag	UNP Q19007
A	4	LEU	-	expression tag	UNP Q19007
A	5	TYR	-	expression tag	UNP Q19007
A	6	PHE	-	expression tag	UNP Q19007
A	7	GLN	-	expression tag	UNP Q19007
A	8	HIS	-	expression tag	UNP Q19007
A	9	MET	-	expression tag	UNP Q19007
B	-11	MET	-	initiating methionine	UNP Q19007
B	-10	GLY	-	expression tag	UNP Q19007
B	-9	SER	-	expression tag	UNP Q19007
B	-8	SER	-	expression tag	UNP Q19007

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Chain	Residue	Modelled	Actual	Comment	Reference
B	-7	HIS	-	expression tag	UNP Q19007
B	-6	HIS	-	expression tag	UNP Q19007
B	-5	HIS	-	expression tag	UNP Q19007
B	-4	HIS	-	expression tag	UNP Q19007
B	-3	HIS	-	expression tag	UNP Q19007
B	-2	HIS	-	expression tag	UNP Q19007
B	-1	SER	-	expression tag	UNP Q19007
B	0	SER	-	expression tag	UNP Q19007
B	1	GLY	-	expression tag	UNP Q19007
B	2	GLU	-	expression tag	UNP Q19007
B	3	ASN	-	expression tag	UNP Q19007
B	4	LEU	-	expression tag	UNP Q19007
B	5	TYR	-	expression tag	UNP Q19007
B	6	PHE	-	expression tag	UNP Q19007
B	7	GLN	-	expression tag	UNP Q19007
B	8	HIS	-	expression tag	UNP Q19007
B	9	MET	-	expression tag	UNP Q19007

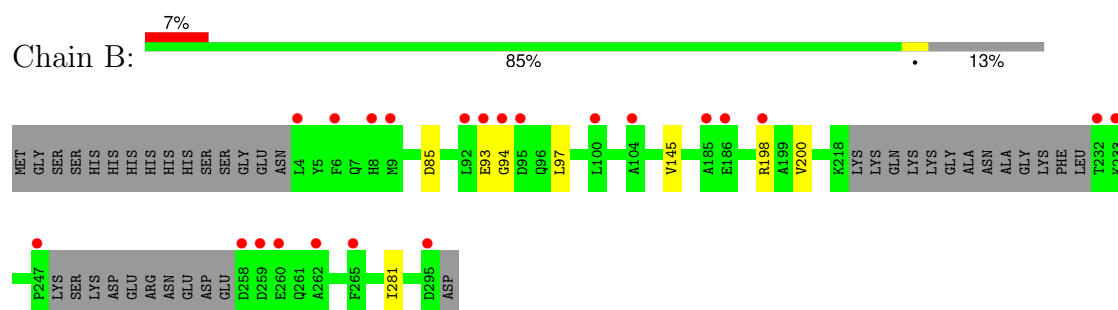
- Molecule 2 is ZINC ION (three-letter code: ZN) (formula: Zn) (labeled as "Ligand of Interest" by depositor).

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

- Molecule 3 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	32	Total O 32 32	0	0
3	B	29	Total O 29 29	0	0

- Molecule 1: Nucleosome Assembly Protein



4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	95.73Å 126.75Å 118.19Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	37.20 – 2.16 37.20 – 2.16	Depositor EDS
% Data completeness (in resolution range)	98.2 (37.20-2.16) 98.2 (37.20-2.16)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.29 (at 2.16Å)	Xtriage
Refinement program	PHENIX (1.14_3260: ???)	Depositor
R, R_{free}	0.250 , 0.279 0.252 , 0.257	Depositor DCC
R_{free} test set	1915 reflections (5.03%)	wwPDB-VP
Wilson B-factor (Å ²)	24.9	Xtriage
Anisotropy	0.118	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.40 , 51.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.90	EDS
Total number of atoms	4299	wwPDB-VP
Average B, all atoms (Å ²)	31.0	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 35.32 % 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 5.9384e-04. 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	0.26	0/2133	0.42	0/2890
1	B	0.26	0/2195	0.42	0/2972
All	All	0.26	0/4328	0.42	0/5862

There are no bond length outliers.

There are no bond angle outliers.

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	2087	0	2029	7	0
1	B	2149	0	2094	4	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	32	0	0	0	0
3	B	29	0	0	0	0
All	All	4299	0	4123	11	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 1.

All (11) 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:B:93:GLU:HG2	1:B:94:GLY:H	1.63	0.63
1:A:38:MET:SD	1:A:188:PRO:HB2	2.47	0.55
1:B:145:VAL:HG12	1:B:281:ILE:HD11	1.95	0.48
1:B:198:ARG:NH2	1:B:200:VAL:CG2	2.77	0.48
1:A:92:LEU:HD12	1:A:97:LEU:HD13	1.97	0.46
1:B:94:GLY:HA2	1:B:97:LEU:HB3	1.98	0.46
1:A:175:LYS:HE2	1:A:243:PHE:O	2.17	0.44
1:A:175:LYS:HD3	1:A:177:TYR:CZ	2.53	0.43
1:A:145:VAL:HG12	1:A:281:ILE:HD11	2.01	0.43
1:A:93:GLU:H	1:A:96:GLN:HB2	1.85	0.42
1:A:52:HIS:CE1	1:A:149:ALA:HB1	2.54	0.41

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	257/308 (83%)	255 (99%)	2 (1%)	0	100	100
1	B	263/308 (85%)	262 (100%)	1 (0%)	0	100	100
All	All	520/616 (84%)	517 (99%)	3 (1%)	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	219/265 (83%)	218 (100%)	1 (0%)	88	92
1	B	230/265 (87%)	229 (100%)	1 (0%)	91	93
All	All	449/530 (85%)	447 (100%)	2 (0%)	91	93

All (2) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	294	SER
1	B	85	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

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 monosaccharides in this entry.

5.6 Ligand geometry [i](#)

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

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers [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	263/308 (85%)	0.36	21 (7%) 12 17	15, 28, 59, 80	0
1	B	269/308 (87%)	0.30	22 (8%) 11 16	14, 28, 57, 97	0
All	All	532/616 (86%)	0.33	43 (8%) 12 16	14, 28, 58, 97	0

All (43) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	93	GLU	5.3
1	B	258	ASP	5.1
1	A	8	HIS	5.0
1	A	235	VAL	4.7
1	A	94	GLY	4.1
1	A	247	PRO	4.1
1	B	232	THR	3.6
1	B	259	ASP	3.6
1	B	295	ASP	3.5
1	B	247	PRO	3.4
1	A	261	GLN	3.4
1	B	4	LEU	3.4
1	B	186	GLU	3.3
1	B	8	HIS	3.3
1	A	4	LEU	3.3
1	A	95	ASP	3.2
1	A	246	PRO	3.1
1	B	233	LYS	3.0
1	A	82	GLU	3.0
1	B	95	ASP	2.8
1	A	186	GLU	2.8
1	A	236	LYS	2.8
1	B	260	GLU	2.8
1	B	265	PHE	2.8

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Mol	Chain	Res	Type	RSRZ
1	B	92	LEU	2.7
1	B	94	GLY	2.7
1	A	3	ASN	2.6
1	A	93	GLU	2.6
1	A	102	LYS	2.5
1	A	234	THR	2.4
1	B	262	ALA	2.4
1	B	100	LEU	2.4
1	A	79	PRO	2.4
1	B	6	PHE	2.3
1	A	84	ILE	2.3
1	A	216	ALA	2.3
1	B	185	ALA	2.2
1	A	215	LYS	2.2
1	B	104	ALA	2.2
1	A	152	ASP	2.1
1	B	9	MET	2.1
1	B	198	ARG	2.0
1	A	6	PHE	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 monosaccharides 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.

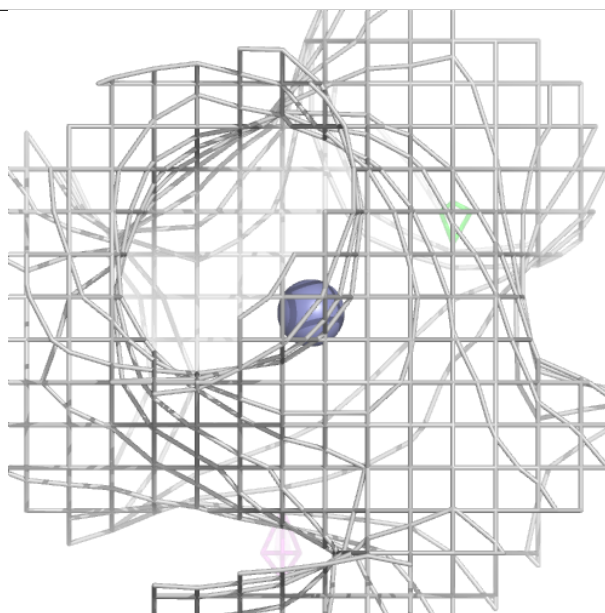
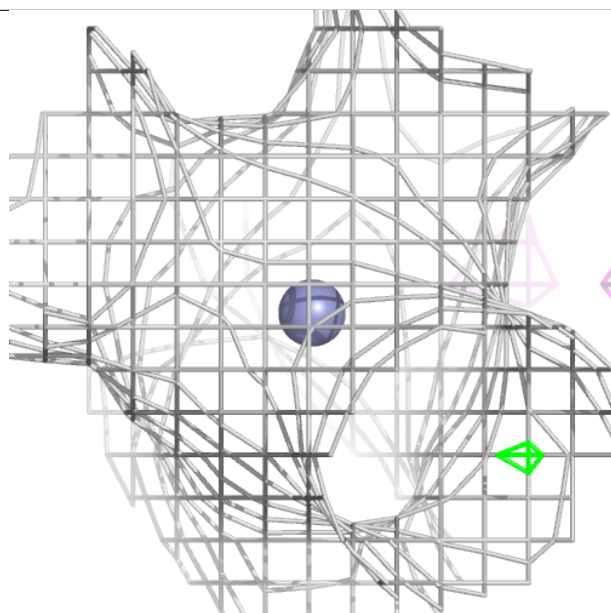
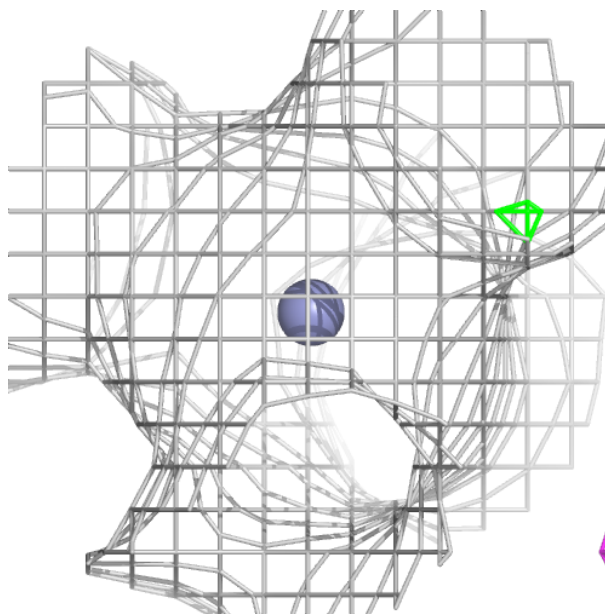
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
2	ZN	A	301	1/1	1.00	0.12	19,19,19,19	0
2	ZN	B	301	1/1	1.00	0.12	19,19,19,19	0

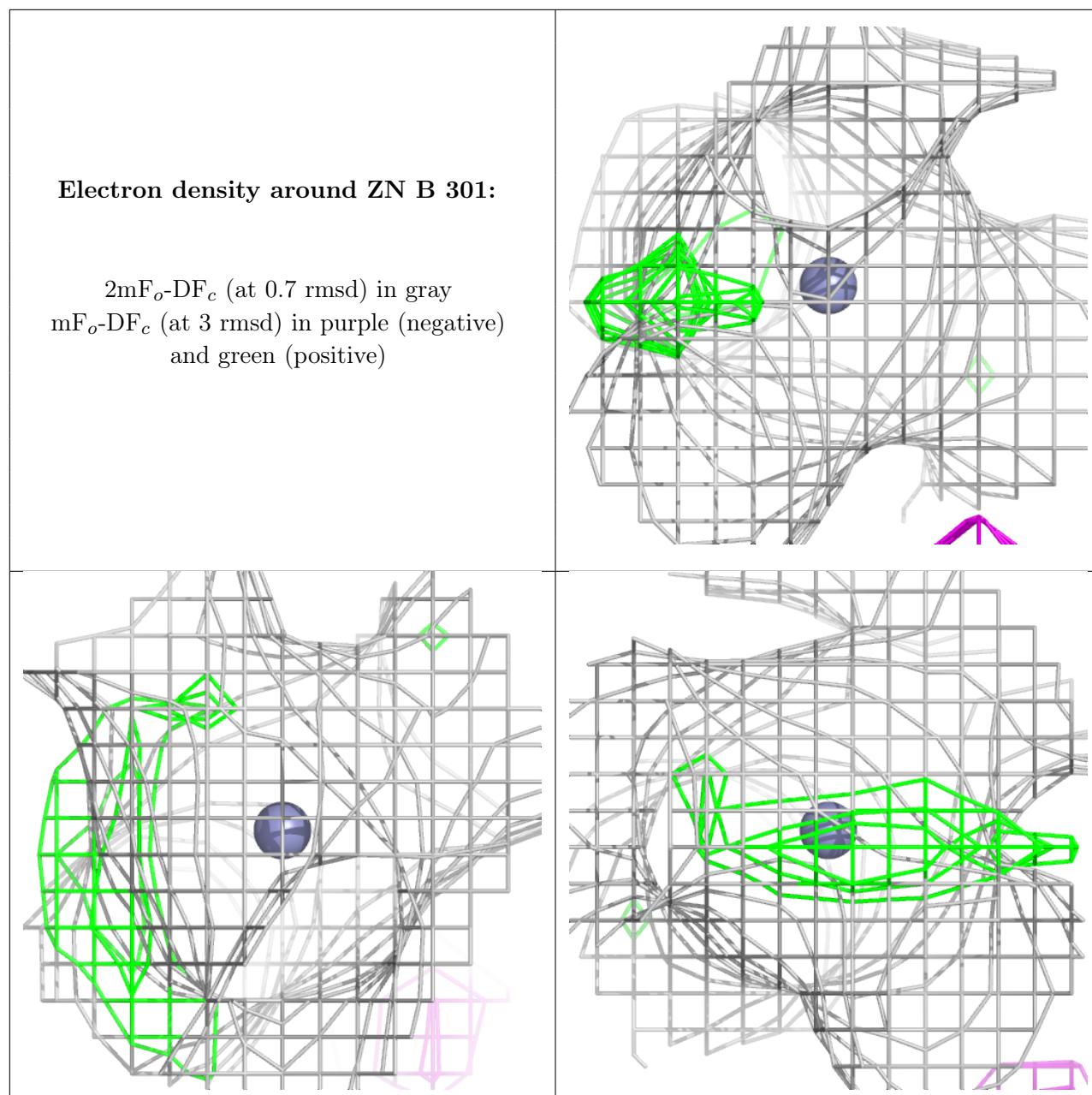
The following is a graphical depiction of the model fit to experimental electron density of all

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 301:

$2mF_o-DF_c$ (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.