

wwPDB X-ray Structure Validation Summary Report (i)

Sep 28, 2021 – 08:26 AM JST

PDB ID 7DBU

> Title Crystal structure of catalytic domain of Anhydrobiosis-related Mn-dependent

> > Peroxidase (AMNP) from Ramazzottius varieornatus (Zn2+-bound form)

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2020-10-21 Deposited on

Resolution 1.60 Å(reported)

This is a wwPDB X-ray Structure Validation Summary 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 (i) symbol.

The following versions of software and data (see references (1)) were used in the production of this report:

MolProbity 4.02b-467Xtriage (Phenix) 1.13

EDS 2.23.2buster-report 1.1.7 (2018)

20191225.v01 (using entries in the PDB archive December 25th 2019) Percentile statistics

> Refmac 5.8.0158

CCP4 7.0.044 (Gargrove) Ideal geometry (proteins) Engh & Huber (2001)

Ideal geometry (DNA, RNA) Parkinson et al. (1996)

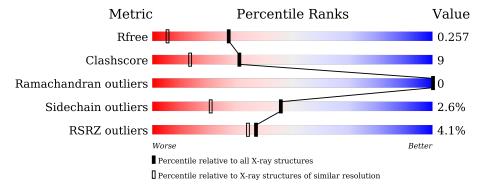
Validation Pipeline (wwPDB-VP) 2.23.2

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure: X- $RAY\ DIFFRACTION$

The reported resolution of this entry is 1.60 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	$\begin{array}{c} \textbf{Whole archive} \\ (\#\text{Entries}) \end{array}$	$\begin{array}{c} {\rm Similar\ resolution} \\ (\#{\rm Entries},\ {\rm resolution\ range}({\rm \AA})) \end{array}$
R_{free}	130704	3398 (1.60-1.60)
Clashscore	141614	3665 (1.60-1.60)
Ramachandran outliers	138981	3564 (1.60-1.60)
Sidechain outliers	138945	3563 (1.60-1.60)
RSRZ outliers	127900	3321 (1.60-1.60)

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 <=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	173	74%	18% 7%
1	В	173	83%	7% • 9%



2 Entry composition (i)

There are 3 unique types of molecules in this entry. The entry contains 2714 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 AMNP/g12777.

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace		
1	Λ	161	Total	С	N	О	S	0	E	0
1	A	101	1246	773	217	248	8	0	9	U
1	D	158	Total	С	N	О	S	0	0	0
1	Б	190	1205	747	213	237	8	0	U	U

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	59	GLY	-	expression tag	UNP A0A1D1VPD8
A	60	SER	-	expression tag	UNP A0A1D1VPD8
A	61	HIS	-	expression tag	UNP A0A1D1VPD8
A	62	MET	-	expression tag	UNP A0A1D1VPD8
В	59	GLY	-	expression tag	UNP A0A1D1VPD8
В	60	SER	-	expression tag	UNP A0A1D1VPD8
В	61	HIS	-	expression tag	UNP A0A1D1VPD8
В	62	MET	-	expression tag	UNP A0A1D1VPD8

• 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	5	Total Zn 5 5	0	0
2	В	6	Total Zn 6 6	0	0

• Molecule 3 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	155	Total O 155 155	0	0

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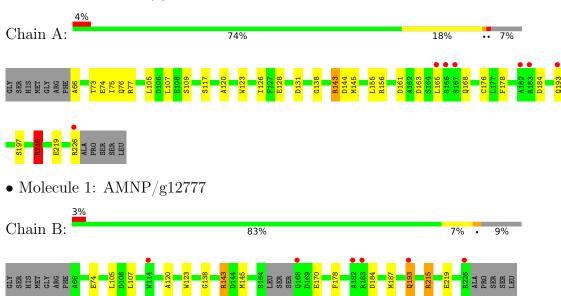
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	В	97	Total O 97 97	0	0



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: AMNP/g12777





4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1	Depositor
Cell constants	41.29Å 48.54Å 49.03Å	Donositor
a, b, c, α , β , γ	73.25° 68.61° 89.95°	Depositor
Resolution (Å)	20.00 - 1.60	Depositor
Resolution (A)	19.66 - 1.60	EDS
% Data completeness	90.0 (20.00-1.60)	Depositor
(in resolution range)	90.1 (19.66-1.60)	EDS
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) > 1$	1.06 (at 1.60Å)	Xtriage
Refinement program	REFMAC 5.8.0253	Depositor
D D.	0.192 , 0.245	Depositor
R, R_{free}	0.201 , 0.257	DCC
R_{free} test set	2014 reflections (5.03%)	wwPDB-VP
Wilson B-factor (Å ²)	9.1	Xtriage
Anisotropy	0.127	Xtriage
Bulk solvent $k_{sol}(e/Å^3)$, $B_{sol}(Å^2)$	0.40, 51.8	EDS
L-test for twinning ²	$ < L >=0.51, < L^2>=0.35$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	2714	wwPDB-VP
Average B, all atoms (Å ²)	19.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: The largest off-origin peak in the Patterson function is 10.16% of the height of the origin peak. No significant pseudotranslation is detected.

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



¹Intensities estimated from amplitudes.

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		nd lengths	Bond angles		
MIOI	Chain	RMSZ # Z > 5		RMSZ	# Z >5	
1	A	0.88	0/1283	1.06	2/1739 (0.1%)	
1	В	0.93	1/1226 (0.1%)	1.04	0/1659	
All	All	0.90	$1/2509 \ (0.0\%)$	1.05	2/3398 (0.1%)	

All (1) bond length outliers are listed below:

\mathbf{Mol}	Chain	Res	Type	Atoms	\mathbf{Z}	$\operatorname{Observed}(\text{\AA})$	$\operatorname{Ideal}(ext{\AA})$
1	В	170	GLU	CD-OE2	-5.37	1.19	1.25

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms Z		$Observed(^o)$	$Ideal(^{o})$
1	A	156	ARG	NE-CZ-NH1	-5.79	117.41	120.30
1	A	215	ARG	CG-CD-NE	-5.05	101.19	111.80

There are no chirality outliers.

There are no planarity outliers.

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	1246	0	1216	31	0
1	В	1205	0	1165	13	0
2	A	5	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	В	6	0	0	0	0
3	A	155	0	0	13	1
3	В	97	0	0	3	1
All	All	2714	0	2381	44	1

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 9.

The worst 5 of 44 close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	$\begin{array}{c} {\rm Interatomic} \\ {\rm distance} \ ({\rm \AA}) \end{array}$	Clash overlap (Å)
1:A:73:THR:HG21	3:A:522:HOH:O	1.53	1.04
1:A:155:LEU:HB2	3:A:522:HOH:O	1.62	0.97
1:A:107:LEU:HD12	1:A:143:ARG:HD2	1.48	0.93
1:A:107:LEU:HD13	1:A:145:MET:CE	1.97	0.92
1:A:128:GLU:OE2	3:A:401:HOH:O	1.90	0.89

All (1) 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:532:HOH:O	3:B:483:HOH:O[1_445]	2.02	0.18

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	Perce	\mathbf{ntiles}
1	A	164/173~(95%)	158 (96%)	6 (4%)	0	100	100
1	В	154/173 (89%)	149 (97%)	5 (3%)	0	100	100
All	All	318/346 (92%)	307 (96%)	11 (4%)	0	100	100



There are no Ramachandran outliers to report.

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	139/143 (97%)	136 (98%)	3 (2%)	52 27	
1	В	131/143 (92%)	127 (97%)	4 (3%)	40 15	
All	All	270/286 (94%)	263 (97%)	7 (3%)	46 21	

5 of 7 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	В	143	ARG
1	В	187	MET
1	В	215	ARG
1	В	193	GLN
1	A	215	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	76	GLN
1	В	160	ASN
1	В	193	GLN

5.3.3 RNA (i)

There are no RNA molecules in this entry.

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

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



5.5 Carbohydrates (i)

There are no monosaccharides in this entry.

5.6 Ligand geometry (i)

Of 11 ligands modelled in this entry, 11 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 (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, 95^{th} 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 $>$	$\# \mathrm{RSRZ}{>}2$	$OWAB(Å^2)$	Q<0.9
1	A	161/173 (93%)	0.49	7 (4%) 35 32	10, 16, 27, 51	0
1	В	158/173 (91%)	0.44	6 (3%) 40 37	11, 16, 31, 59	0
All	All	319/346 (92%)	0.46	13 (4%) 37 34	10, 16, 30, 59	0

The worst 5 of 13 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	166	SER	4.5
1	В	226	ARG	4.2
1	В	183	ARG	4.0
1	A	183	ARG	4.0
1	В	182	ALA	3.6

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)

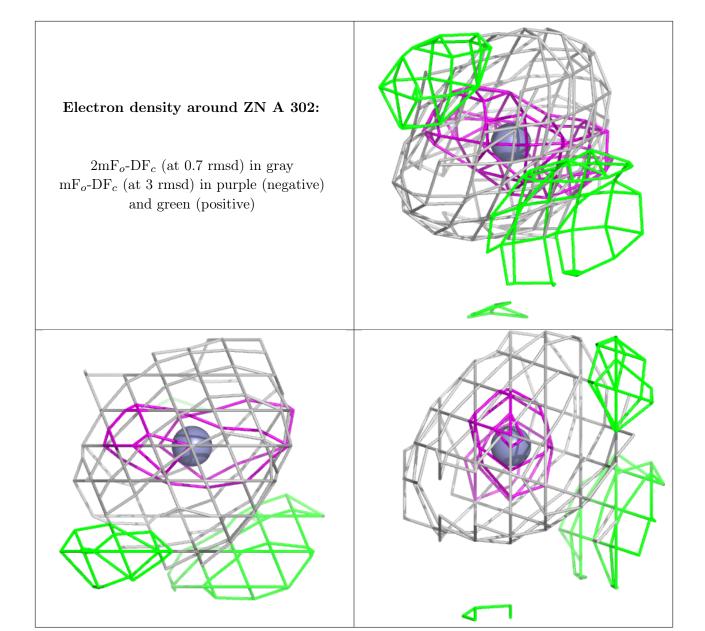
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95^{th} percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9'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



Mol	Type	Chain	Res	Atoms	RSCC	RSR	$\mathbf{B} ext{-}\mathbf{factors}(\mathbf{\mathring{A}}^2)$	Q<0.9
2	ZN	A	302	1/1	0.99	0.13	31,31,31,31	0
2	ZN	A	303	1/1	0.99	0.11	36,36,36,36	0
2	ZN	В	304	1/1	0.99	0.18	38,38,38,38	0
2	ZN	В	305	1/1	0.99	0.13	36,36,36,36	0
2	ZN	A	305	1/1	1.00	0.03	23,23,23,23	1
2	ZN	В	301	1/1	1.00	0.04	13,13,13,13	0
2	ZN	В	302	1/1	1.00	0.05	18,18,18,18	0
2	ZN	В	303	1/1	1.00	0.03	13,13,13,13	0
2	ZN	A	301	1/1	1.00	0.04	18,18,18,18	0
2	ZN	A	304	1/1	1.00	0.11	37,37,37,37	0
2	ZN	В	306	1/1	1.00	0.04	20,20,20,20	1

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.





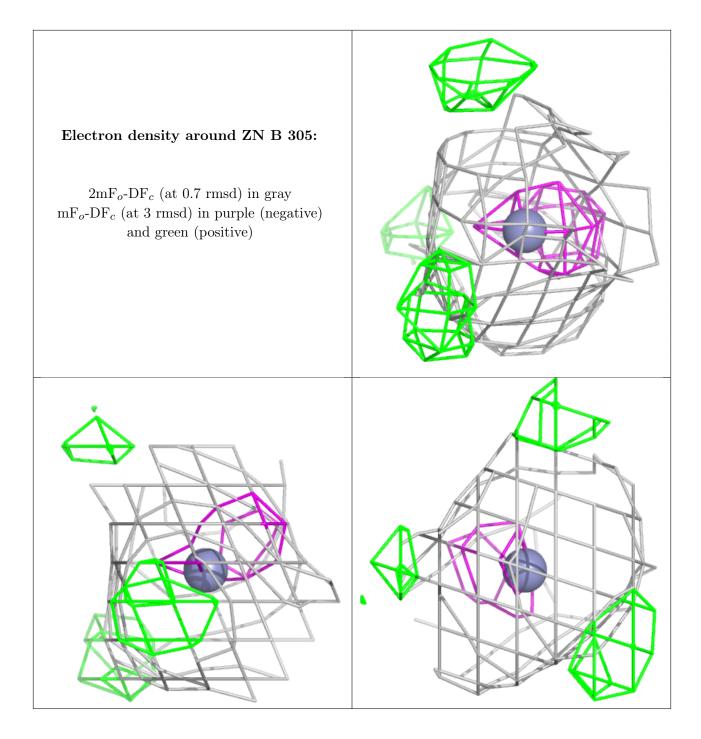


Electron density around ZN A 303: $2 \mathrm{mF}_o\text{-}\mathrm{DF}_c$ (at 0.7 rmsd) in gray $\mathrm{mF}_o\text{-}\mathrm{DF}_c$ (at 3 rmsd) in purple (negative) and green (positive)

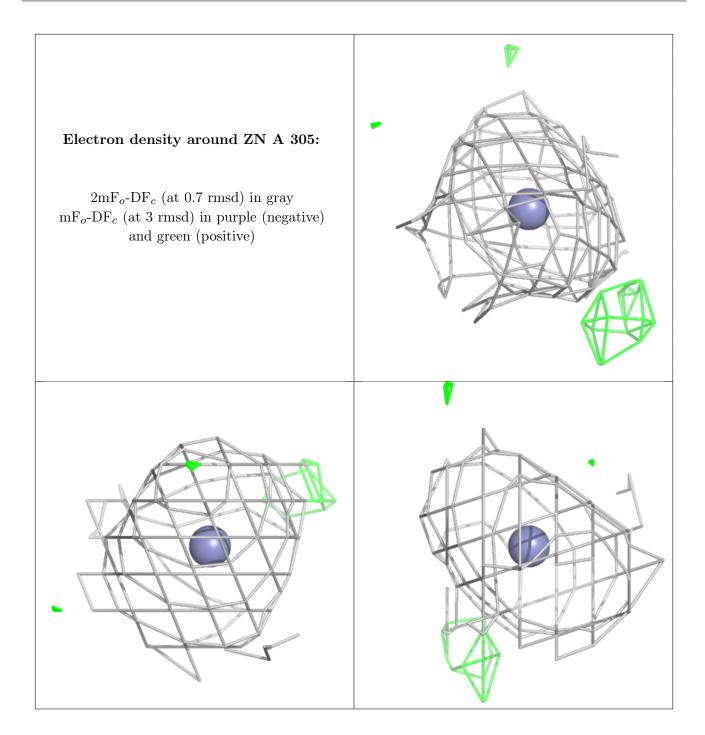


Electron density around ZN B 304: $2 \mathrm{mF}_o\text{-}\mathrm{DF}_c$ (at 0.7 rmsd) in gray $\mathrm{mF}_o\text{-}\mathrm{DF}_c$ (at 3 rmsd) in purple (negative) and green (positive)

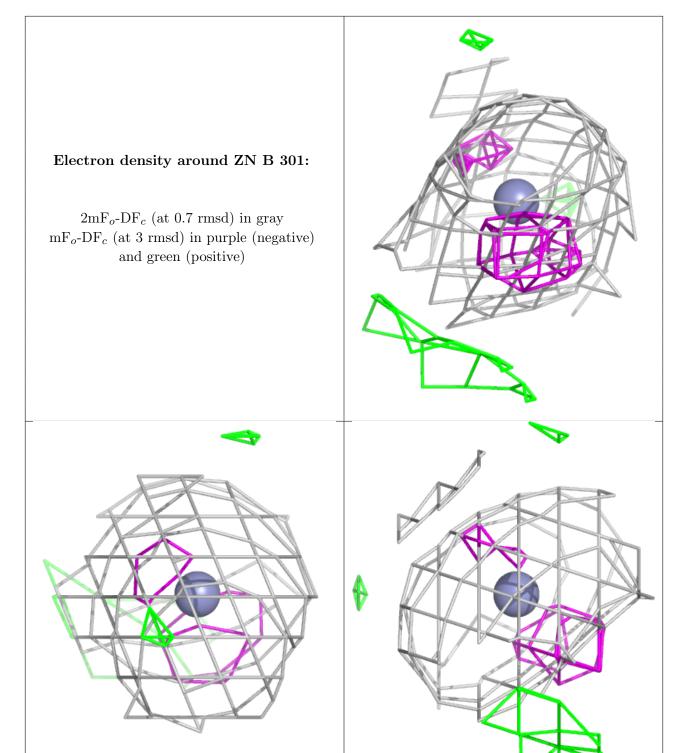








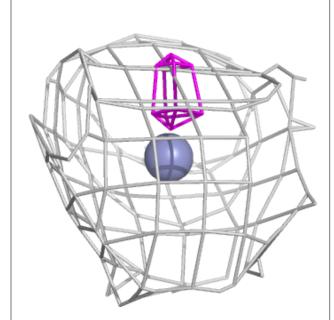


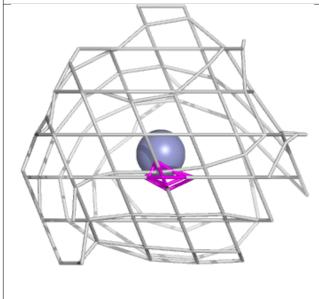


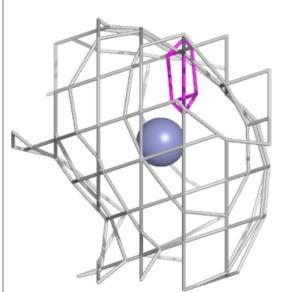


Electron density around ZN B 302:

 $2 {
m mF}_o {
m -DF}_c$ (at 0.7 rmsd) in gray ${
m mF}_o {
m -DF}_c$ (at 3 rmsd) in purple (negative) and green (positive)

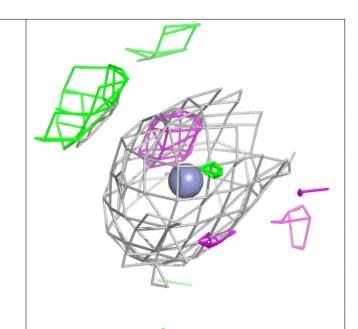


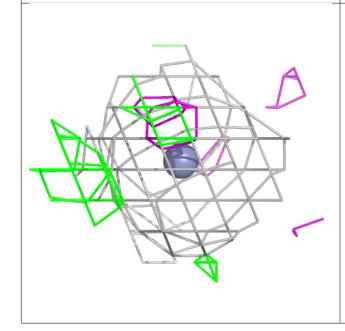


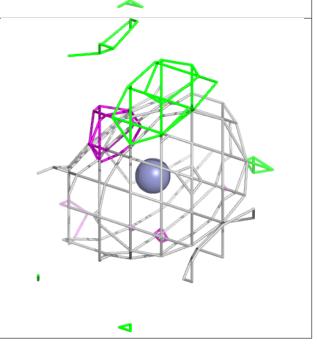


Electron density around ZN B 303:

 $2 {
m mF}_o {
m -DF}_c$ (at 0.7 rmsd) in gray ${
m mF}_o {
m -DF}_c$ (at 3 rmsd) in purple (negative) and green (positive)



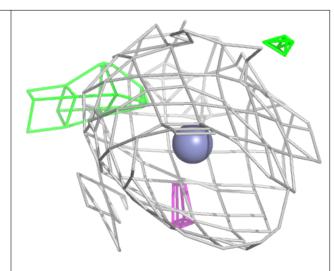


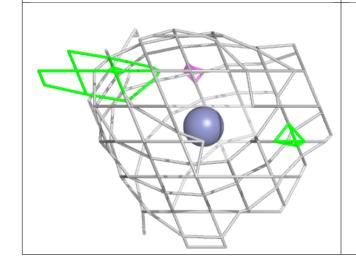


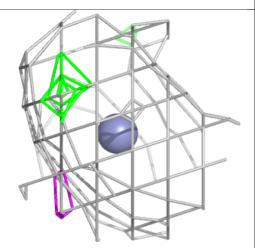


Electron density around ZN A 301:

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



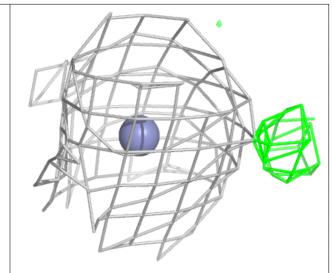


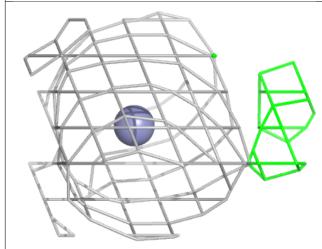


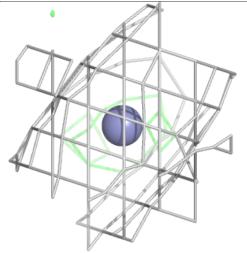


Electron density around ZN A 304:

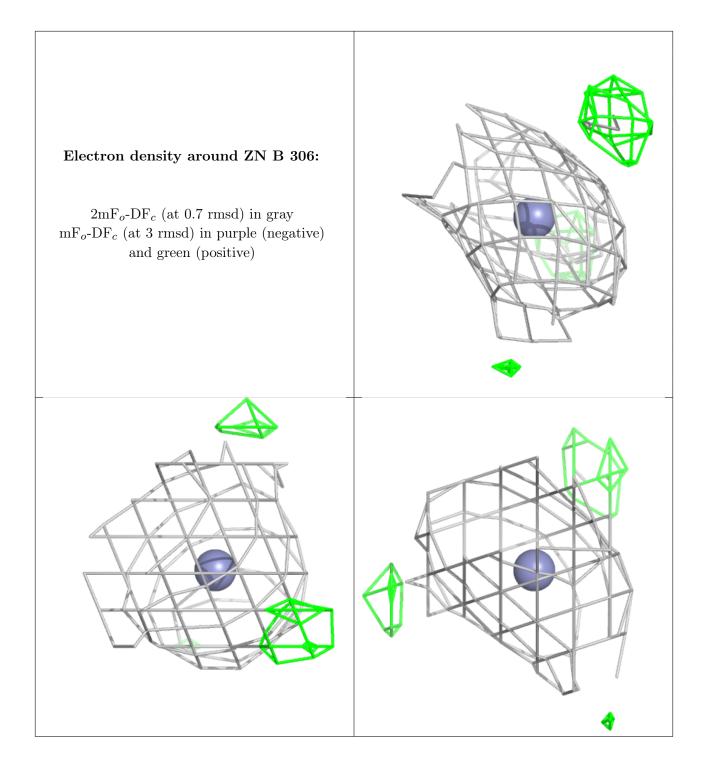
 $2 {
m mF}_o {
m -DF}_c$ (at 0.7 rmsd) in gray ${
m mF}_o {
m -DF}_c$ (at 3 rmsd) in purple (negative) and green (positive)











6.5 Other polymers (i)

There are no such residues in this entry.

