



Full wwPDB X-ray Structure Validation Report ⓘ

Sep 10, 2020 – 09:36 AM BST

PDB ID : 6T5L
Title : MYO-1 from Myroides odoratimimus. Environmental metallo-beta-lactamase
s exhibit high enzymatic activity under zinc deprivation
Authors : Frohlich, C.
Deposited on : 2019-10-16
Resolution : 2.17 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 following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.14.3.dev2
buster-report	:	1.1.7 (2018)
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)	:	2.14.3.dev2

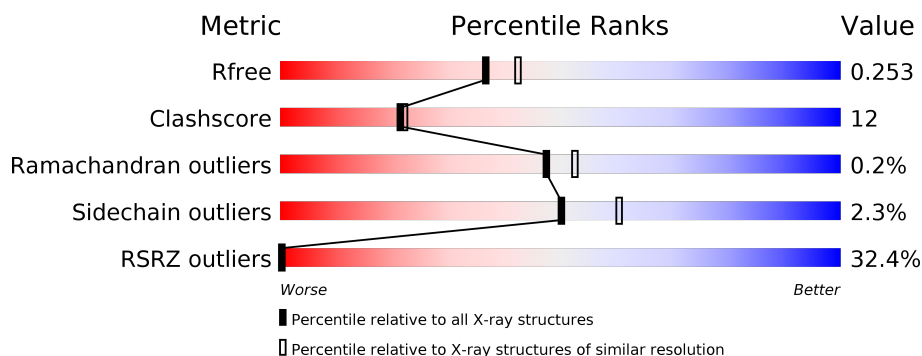
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.17 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	6864 (2.20-2.16)
Clashscore	141614	7689 (2.20-2.16)
Ramachandran outliers	138981	7564 (2.20-2.16)
Sidechain outliers	138945	7564 (2.20-2.16)
RSRZ outliers	127900	6738 (2.20-2.16)

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 on 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	238	<div> <div>18%</div> <div> <div></div> <div>83%</div> <div>11%</div> <div>6%</div> </div> </div>
1	B	238	<div> <div>41%</div> <div> <div></div> <div>59%</div> <div>27%</div> <div>13%</div> </div> </div>

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 6963 atoms, of which 3358 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 Subclass B1 metallo-beta-lactamase.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	224	Total	C	H	N	O	S	57	3	0
			3541	1150	1746	285	356	4			
1	B	208	Total	C	H	N	O	S	350	2	0
			3275	1067	1612	265	327	4			

There are 18 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	18	GLY	-	expression tag	UNP A0A0U3H0V9
A	19	GLN	-	expression tag	UNP A0A0U3H0V9
A	20	GLU	-	expression tag	UNP A0A0U3H0V9
A	21	ASN	-	expression tag	UNP A0A0U3H0V9
A	22	LYS	-	expression tag	UNP A0A0U3H0V9
A	23	LYS	-	expression tag	UNP A0A0U3H0V9
A	24	GLU	-	expression tag	UNP A0A0U3H0V9
A	25	ILE	-	expression tag	UNP A0A0U3H0V9
A	26	ILE	-	expression tag	UNP A0A0U3H0V9
B	18	GLY	-	expression tag	UNP A0A0U3H0V9
B	19	GLN	-	expression tag	UNP A0A0U3H0V9
B	20	GLU	-	expression tag	UNP A0A0U3H0V9
B	21	ASN	-	expression tag	UNP A0A0U3H0V9
B	22	LYS	-	expression tag	UNP A0A0U3H0V9
B	23	LYS	-	expression tag	UNP A0A0U3H0V9
B	24	GLU	-	expression tag	UNP A0A0U3H0V9
B	25	ILE	-	expression tag	UNP A0A0U3H0V9
B	26	ILE	-	expression tag	UNP A0A0U3H0V9

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

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

- Molecule 3 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total 1	Mg 1	0	0

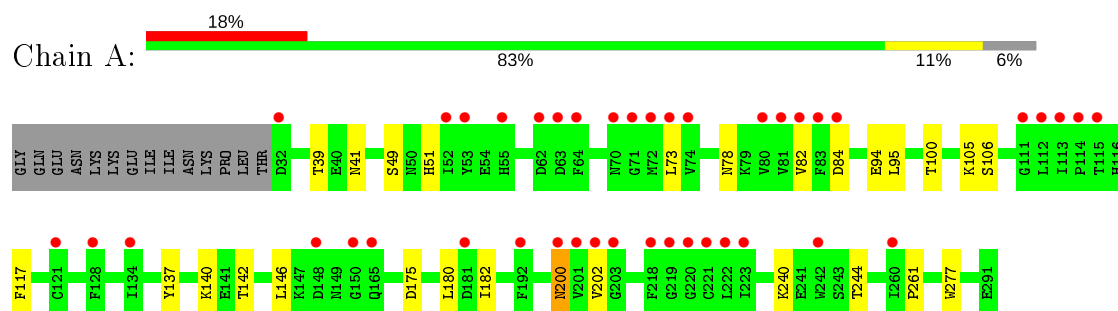
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	114	Total 114	O 114	0	0
4	B	28	Total 28	O 28	0	0

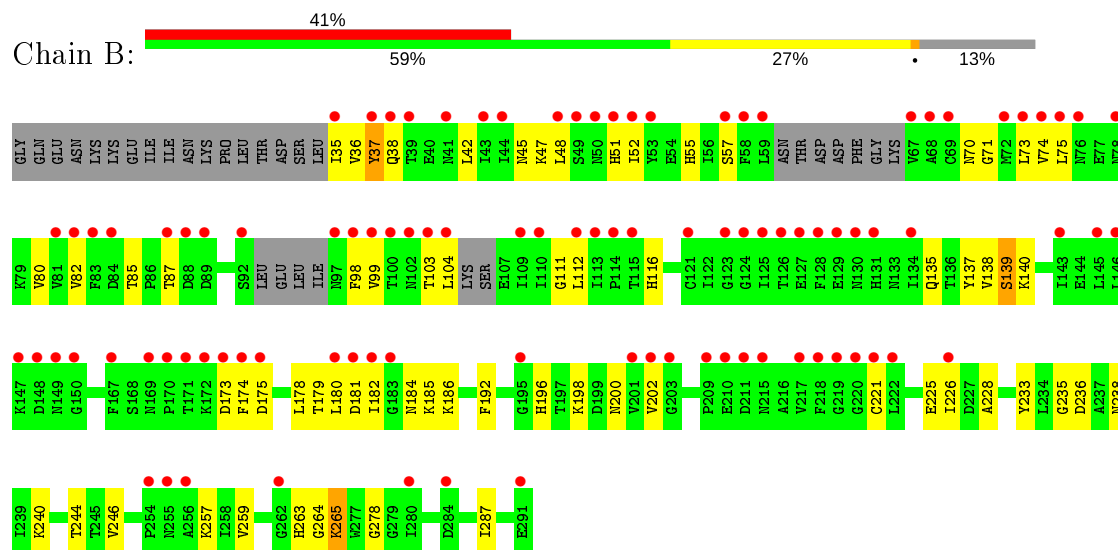
3 Residue-property plots [i](#)

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: Subclass B1 metallo-beta-lactamase



- Molecule 1: Subclass B1 metallo-beta-lactamase



4 Data and refinement statistics

Property	Value	Source
Space group	P 65	Depositor
Cell constants a, b, c, α , β , γ	144.68 Å 144.68 Å 53.31 Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	24.53 – 2.17 24.53 – 2.17	Depositor EDS
% Data completeness (in resolution range)	98.8 (24.53-2.17) 98.8 (24.53-2.17)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.31 (at 2.17 Å)	Xtriage
Refinement program	PHENIX 1.14 _3260	Depositor
R, R_{free}	0.217 , 0.253 0.217 , 0.253	Depositor DCC
R_{free} test set	1655 reflections (4.93%)	wwPDB-VP
Wilson B-factor (Å ²)	48.9	Xtriage
Anisotropy	0.091	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 55.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.35$	Xtriage
Estimated twinning fraction	0.032 for h,-h-k,-l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	6963	wwPDB-VP
Average B, all atoms (Å ²)	74.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.34% 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, MG

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	0.75	1/1837 (0.1%)	0.80	0/2488
1	B	0.53	0/1704	0.69	0/2305
All	All	0.65	1/3541 (0.0%)	0.75	0/4793

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	49	SER	CB-OG	5.15	1.49	1.42

There are no bond angle outliers.

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	1795	1746	1742	19	0
1	B	1663	1612	1602	57	0
2	A	2	0	0	1	0
2	B	2	0	0	0	0
3	A	1	0	0	0	0
4	A	114	0	0	5	3
4	B	28	0	0	6	0
All	All	3605	3358	3344	77	3

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 12.

All (77) 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:142:THR:OG1	4:A:401:HOH:O	1.74	1.03
1:A:73:LEU:HD22	1:A:82:VAL:HG22	1.55	0.87
1:B:175:ASP:O	4:B:401:HOH:O	1.95	0.83
2:A:301:ZN:ZN	4:A:416:HOH:O	1.26	0.82
1:A:39:THR:HG22	1:A:41:ASN:H	1.45	0.81
1:A:51:HIS:NE2	4:A:402:HOH:O	2.16	0.78
1:B:55:HIS:ND1	1:B:73:LEU:HD12	2.00	0.77
1:B:36:VAL:O	1:B:36:VAL:HG12	1.84	0.76
1:B:35:ILE:HG12	1:B:36:VAL:HG23	1.68	0.75
1:B:70:ASN:O	4:B:402:HOH:O	2.05	0.75
1:A:73:LEU:CD2	1:A:82:VAL:HG22	2.15	0.75
1:B:52:ILE:CD1	1:B:74:VAL:HG22	2.17	0.73
1:B:236:ASP:OD1	4:B:403:HOH:O	2.08	0.70
1:A:73:LEU:HD11	1:A:95:LEU:HD11	1.75	0.68
1:A:39:THR:HG21	1:A:94:GLU:OE1	1.95	0.67
1:B:55:HIS:ND1	1:B:73:LEU:CD1	2.57	0.67
1:B:52:ILE:HD12	1:B:74:VAL:HG22	1.77	0.65
1:B:51:HIS:HB3	1:B:75:LEU:O	1.96	0.65
1:B:55:HIS:NE2	1:B:71:GLY:HA3	2.12	0.63
1:B:55:HIS:HE1	1:B:82:VAL:CG2	2.12	0.63
1:B:99:VAL:O	1:B:104:LEU:N	2.31	0.63
1:B:75:LEU:HB2	1:B:80:VAL:HG22	1.82	0.62
1:B:182:ILE:N	1:B:185:LYS:O	2.33	0.60
1:B:244:THR:HG23	4:B:421:HOH:O	2.01	0.60
1:A:244:THR:HG23	4:A:421:HOH:O	2.02	0.58
1:B:55:HIS:CE1	1:B:82:VAL:CG2	2.89	0.56
1:B:35:ILE:O	1:B:45:ASN:HA	2.07	0.55
1:B:47:LYS:C	1:B:48:LEU:HD23	2.28	0.54
1:B:265:LYS:CD	1:B:265:LYS:H	2.22	0.53
1:A:182:ILE:HG23	1:A:182:ILE:O	2.09	0.53
1:B:225:GLU:HG2	1:B:228:ALA:HB2	1.91	0.53
1:A:244:THR:HG22	4:A:422:HOH:O	2.07	0.53
1:B:137:TYR:CD1	1:B:180:LEU:HD11	2.44	0.52
1:B:198:LYS:H	1:B:198:LYS:HE2	1.75	0.52
1:B:116:HIS:HE1	1:B:221:CYS:SG	2.32	0.52
1:A:117:PHE:HB2	1:A:146:LEU:HG	1.91	0.51
1:B:98:PHE:O	1:B:103:THR:HB	2.11	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:200[B]:ASN:C	1:A:200[B]:ASN:HD22	2.15	0.50
1:B:35:ILE:HG12	1:B:36:VAL:CG2	2.40	0.50
1:B:35:ILE:HG23	1:B:36:VAL:HG23	1.92	0.50
1:B:198:LYS:CE	1:B:198:LYS:H	2.25	0.49
1:A:137:TYR:CZ	1:A:180:LEU:HD21	2.49	0.48
1:B:140:LYS:HG2	1:B:173:ASP:HB3	1.96	0.48
1:A:140:LYS:HD2	1:A:175:ASP:OD1	2.13	0.47
1:B:178:LEU:HD11	1:B:180:LEU:HG	1.96	0.47
1:B:265:LYS:HD2	1:B:265:LYS:H	1.78	0.47
1:B:259:VAL:N	1:B:278:GLY:O	2.44	0.47
1:B:265:LYS:HD3	4:B:410:HOH:O	2.14	0.47
1:B:36:VAL:O	1:B:37:TYR:C	2.53	0.47
1:B:137:TYR:CE1	1:B:180:LEU:HD11	2.51	0.46
1:B:238:ASN:OD1	1:B:240:LYS:HG3	2.16	0.45
1:B:36:VAL:O	1:B:36:VAL:CG1	2.54	0.45
1:B:111:GLY:HA2	1:B:135:GLN:O	2.17	0.44
1:B:57:SER:HB2	1:B:87:THR:H	1.82	0.44
1:B:178:LEU:HD12	1:B:179:THR:N	2.32	0.44
1:A:100:THR:O	1:A:105:LYS:HA	2.17	0.44
1:B:181:ASP:OD2	1:B:186:LYS:HE3	2.18	0.43
1:B:192:PHE:HE1	1:B:202:VAL:HG22	1.84	0.43
1:B:225:GLU:OE2	1:B:264:GLY:HA3	2.19	0.43
1:B:225:GLU:OE2	1:B:265:LYS:HD2	2.19	0.43
1:B:246:VAL:CG1	1:B:287:ILE:HG13	2.49	0.43
1:A:78:ASN:HA	1:A:106:SER:OG	2.19	0.42
1:A:261:PRO:HD3	1:A:277:TRP:HA	2.02	0.42
1:B:139:SER:HA	1:B:174:PHE:O	2.19	0.42
1:B:263:HIS:O	4:B:405:HOH:O	2.21	0.42
1:B:196:HIS:CD2	1:B:221:CYS:HB2	2.55	0.41
1:B:226:ILE:HG22	1:B:265:LYS:NZ	2.35	0.41
1:A:240:LYS:HA	1:A:240:LYS:HD2	1.83	0.41
1:A:137:TYR:CE2	1:A:180:LEU:HD21	2.55	0.41
1:B:99:VAL:O	1:B:104:LEU:CA	2.68	0.41
1:B:198:LYS:NZ	1:B:236:ASP:HB3	2.36	0.41
1:B:226:ILE:HA	1:B:226:ILE:HD12	1.89	0.41
1:B:85:THR:CG2	1:B:112:LEU:HD21	2.51	0.41
1:B:198:LYS:HB2	1:B:198:LYS:HE2	1.83	0.40
1:B:36:VAL:O	1:B:38:GLN:N	2.53	0.40
1:B:233:TYR:CZ	1:B:235:GLY:HA3	2.55	0.40
1:B:138:VAL:O	1:B:173:ASP:HA	2.22	0.40

All (3) symmetry-related close contacts are listed below. The label for Atom-2 includes the sym-

metry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:A:402:HOH:O	4:A:455:HOH:O[3_655]	1.99	0.21
4:A:436:HOH:O	4:A:489:HOH:O[3_655]	2.13	0.07
4:A:455:HOH:O	4:A:480:HOH:O[2_544]	2.18	0.02

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	225/238 (94%)	218 (97%)	6 (3%)	1 (0%)	34	35
1	B	202/238 (85%)	190 (94%)	12 (6%)	0	100	100
All	All	427/476 (90%)	408 (96%)	18 (4%)	1 (0%)	47	52

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	84	ASP

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	202/213 (95%)	198 (98%)	4 (2%)	55	66
1	B	187/213 (88%)	179 (96%)	8 (4%)	29	34
All	All	389/426 (91%)	377 (97%)	12 (3%)	50	48

All (12) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	200[A]	ASN
1	A	200[B]	ASN
1	A	202[A]	VAL
1	A	202[B]	VAL
1	B	37	TYR
1	B	42	LEU
1	B	139	SER
1	B	184	ASN
1	B	200[A]	ASN
1	B	200[B]	ASN
1	B	257	LYS
1	B	265	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	131	HIS
1	B	149	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 monosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 5 ligands modelled in this entry, 5 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	224/238 (94%)	1.05	42 (18%) 1 1	35, 49, 74, 96	8 (3%)
1	B	208/238 (87%)	2.66	98 (47%) 0 0	47, 85, 118, 162	54 (25%)
All	All	432/476 (90%)	1.82	140 (32%) 0 0	35, 61, 112, 162	62 (14%)

All (140) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	82	VAL	18.3
1	B	92	SER	18.1
1	B	73	LEU	14.8
1	B	53	TYR	12.2
1	B	97	ASN	11.3
1	B	74	VAL	10.2
1	B	98	PHE	10.2
1	B	125	ILE	9.7
1	B	103	THR	9.5
1	B	100	THR	8.9
1	B	99	VAL	8.7
1	B	102	ASN	8.6
1	B	35	ILE	8.5
1	B	150	GLY	8.4
1	B	83	PHE	8.2
1	B	52	ILE	7.8
1	B	218	PHE	6.9
1	B	219	GLY	6.8
1	B	113	ILE	6.8
1	B	202	VAL	6.7
1	B	134	ILE	6.6
1	B	126	THR	6.5
1	B	89	ASP	6.4
1	A	83	PHE	6.3

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Mol	Chain	Res	Type	RSRZ
1	B	58	PHE	6.2
1	A	63	ASP	6.1
1	A	218	PHE	6.0
1	A	72	MET	5.8
1	B	255	ASN	5.5
1	B	201[A]	VAL	5.3
1	A	202[A]	VAL	5.3
1	A	82	VAL	5.2
1	B	203	GLY	5.2
1	B	131	HIS	5.2
1	B	115	THR	5.1
1	B	130	ASN	5.1
1	B	182	ILE	5.0
1	B	124	GLY	5.0
1	B	57	SER	4.9
1	B	149	ASN	4.9
1	B	143	ILE	4.9
1	A	150	GLY	4.8
1	B	38	GLN	4.7
1	B	170	PRO	4.7
1	B	210	GLU	4.7
1	B	147	LYS	4.7
1	A	81	VAL	4.6
1	A	113	ILE	4.6
1	B	104	LEU	4.6
1	A	73	LEU	4.6
1	B	183	GLY	4.5
1	A	219	GLY	4.5
1	B	280	ILE	4.4
1	A	201	VAL	4.3
1	B	87	THR	4.3
1	A	115	THR	4.2
1	B	291	GLU	4.2
1	A	71	GLY	4.1
1	B	68	ALA	4.1
1	A	222	LEU	4.0
1	B	256	ALA	3.8
1	B	220	GLY	3.8
1	A	62	ASP	3.8
1	A	221	CYS	3.8
1	B	44	ILE	3.8
1	B	180	LEU	3.8

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Mol	Chain	Res	Type	RSRZ
1	B	254	PRO	3.7
1	A	260	ILE	3.6
1	B	88	ASP	3.6
1	B	78	ASN	3.6
1	B	112	LEU	3.6
1	A	220	GLY	3.5
1	B	114	PRO	3.4
1	B	39	THR	3.4
1	B	41	ASN	3.4
1	B	67	VAL	3.3
1	B	81	VAL	3.3
1	B	48	LEU	3.3
1	B	172	LYS	3.3
1	A	121	CYS	3.3
1	B	51	HIS	3.2
1	A	112	LEU	3.2
1	A	203[A]	GLY	3.2
1	B	209	PRO	3.2
1	B	226	ILE	3.1
1	B	148	ASP	3.1
1	A	80	VAL	3.1
1	B	50	ASN	3.0
1	B	171	THR	3.0
1	B	84	ASP	3.0
1	B	129	GLU	3.0
1	A	64	PHE	3.0
1	A	128	PHE	2.9
1	A	223	ILE	2.9
1	B	43	ILE	2.9
1	B	128	PHE	2.9
1	B	173	ASP	2.9
1	B	49	SER	2.8
1	B	181	ASP	2.8
1	A	134	ILE	2.8
1	A	111	GLY	2.8
1	B	69	CYS	2.8
1	A	55	HIS	2.7
1	A	148	ASP	2.7
1	B	222	LEU	2.7
1	B	121	CYS	2.6
1	B	167	PHE	2.6
1	B	37	TYR	2.6

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Mol	Chain	Res	Type	RSRZ
1	A	52	ILE	2.6
1	B	169	ASN	2.5
1	B	123	GLY	2.5
1	A	84	ASP	2.5
1	B	72	MET	2.5
1	A	181	ASP	2.5
1	B	217	VAL	2.5
1	B	109	ILE	2.5
1	A	192	PHE	2.4
1	B	146	LEU	2.4
1	B	145	LEU	2.3
1	A	32	ASP	2.3
1	A	53	TYR	2.3
1	B	174	PHE	2.3
1	A	242	TRP	2.3
1	B	211	ASP	2.3
1	B	59	LEU	2.2
1	B	75	LEU	2.2
1	A	74	VAL	2.2
1	B	284	ASP	2.2
1	A	114	PRO	2.2
1	B	195	GLY	2.2
1	B	175	ASP	2.2
1	A	200[A]	ASN	2.2
1	B	215	ASN	2.1
1	B	262	GLY	2.1
1	A	70	ASN	2.1
1	B	110	ILE	2.1
1	B	76	ASN	2.1
1	B	127	GLU	2.1
1	A	165	GLN	2.0
1	B	221	CYS	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

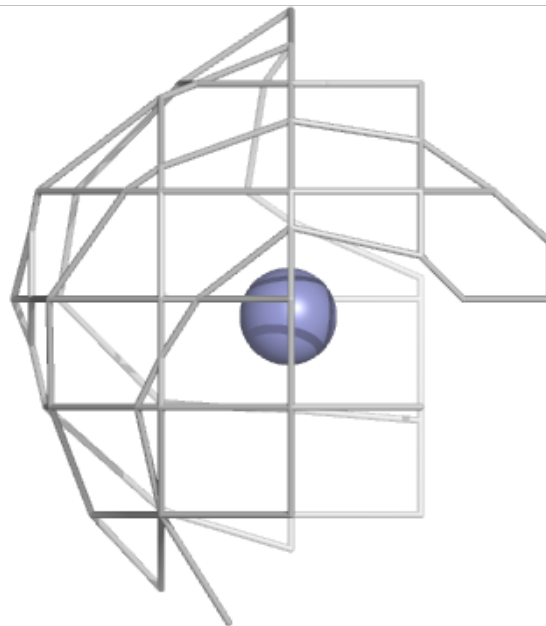
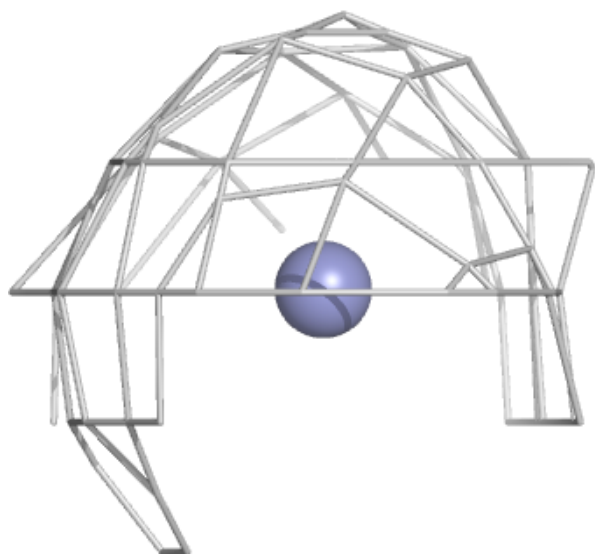
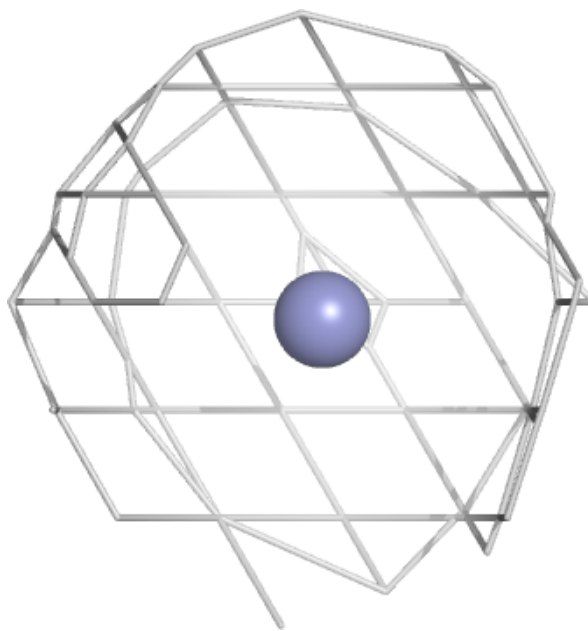
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	ZN	B	302	1/1	0.91	0.07	79,79,79,79	1
3	MG	A	303	1/1	0.96	0.12	47,47,47,47	0
2	ZN	A	302	1/1	0.97	0.12	49,49,49,49	1
2	ZN	A	301	1/1	0.99	0.08	49,49,49,49	0
2	ZN	B	301	1/1	0.99	0.08	70,70,70,70	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.

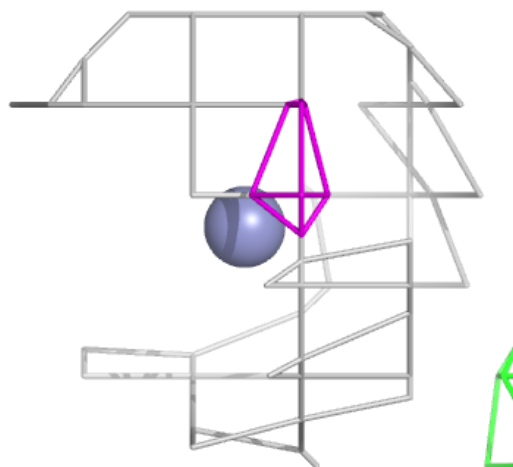
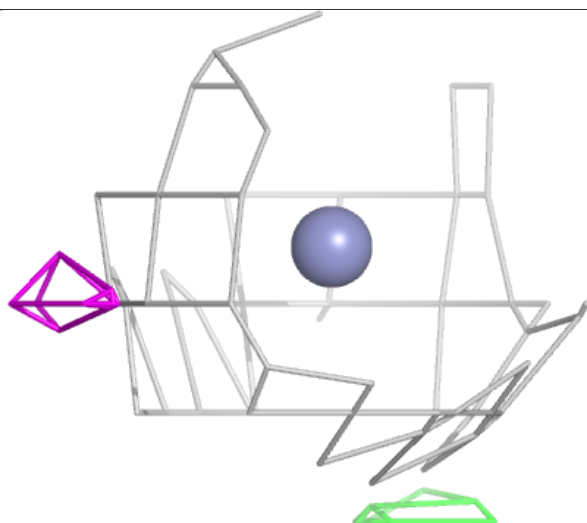
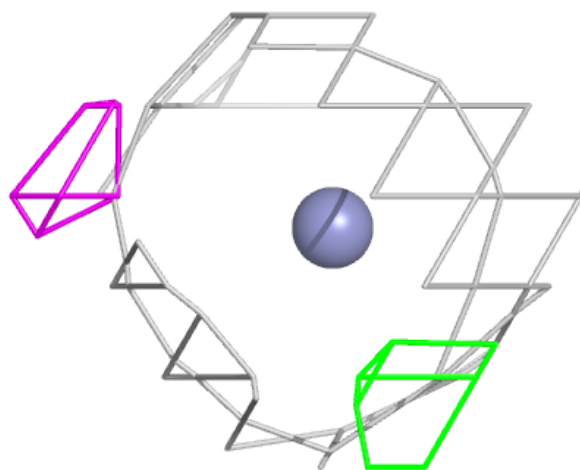
Electron density around ZN B 302:

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



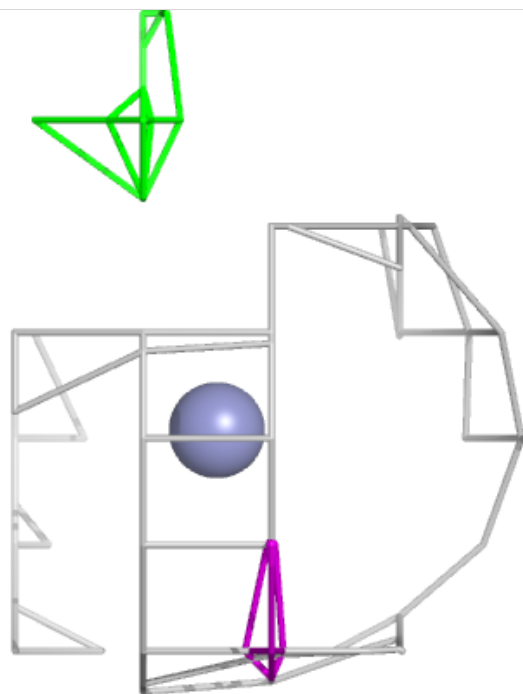
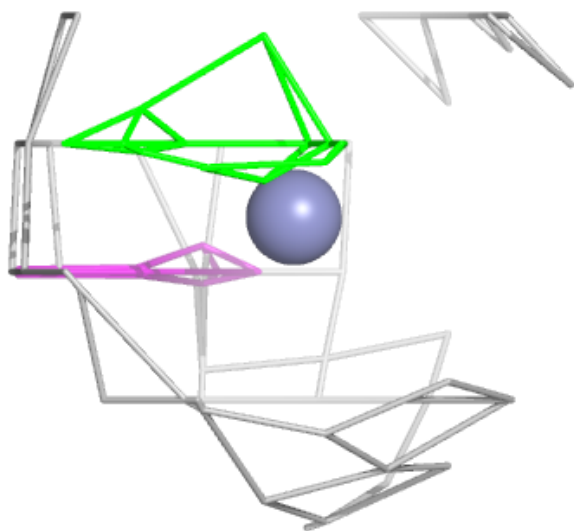
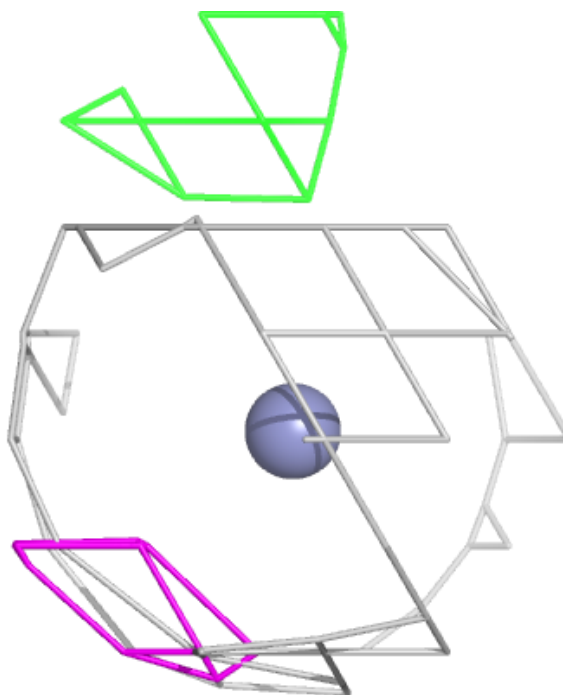
Electron density around ZN A 302:

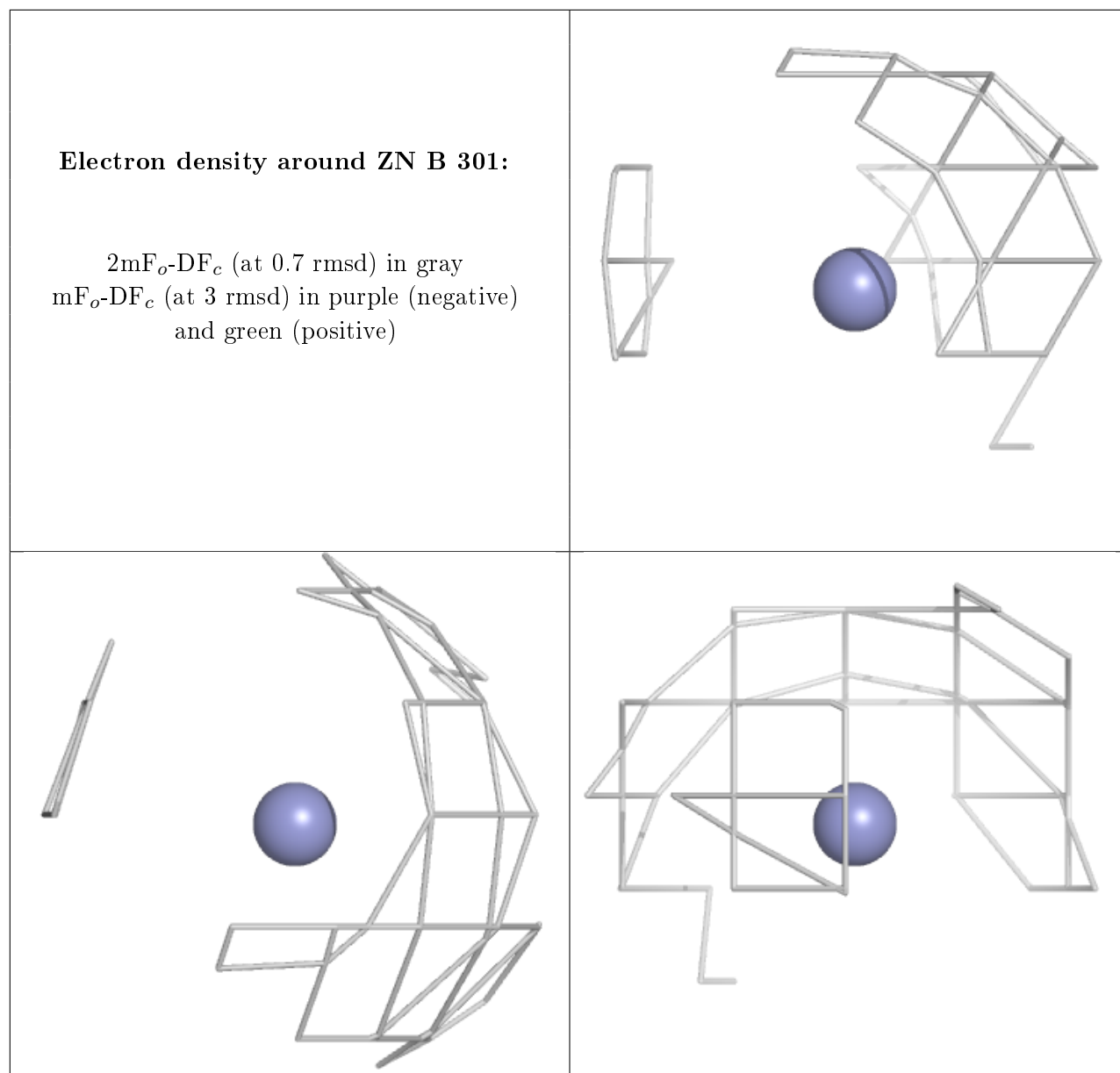
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around ZN A 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.