



wwPDB X-ray Structure Validation Summary Report ⓘ

Oct 18, 2020 – 12:14 AM JST

PDB ID : 6M53
Title : Crystal structure of 2, 3-dihydroxybenzoic acid decarboxylase from *Fusarium oxysporum*
Authors : Song, M.K.; Feng, J.H.; Liu, W.D.; Wu, Q.Q.; Zhu, D.M.
Deposited on : 2020-03-09
Resolution : 1.55 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 ⓘ symbol.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.14.6
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) : 2.14.6

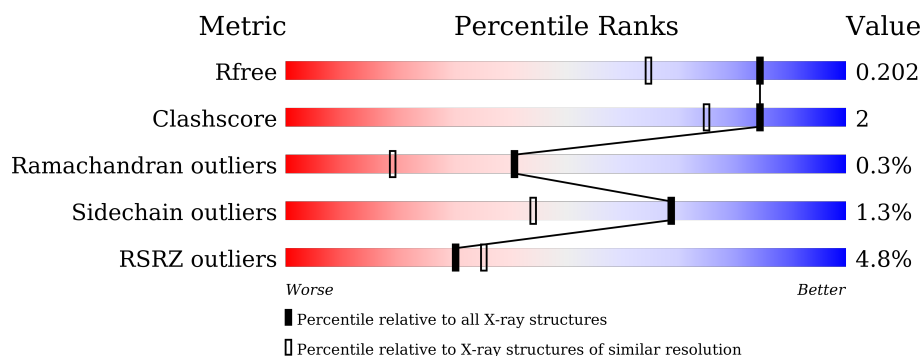
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.55 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	1483 (1.56-1.56)
Clashscore	141614	1529 (1.56-1.56)
Ramachandran outliers	138981	1498 (1.56-1.56)
Sidechain outliers	138945	1495 (1.56-1.56)
RSRZ outliers	127900	1465 (1.56-1.56)

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	343	<div> <div>0%</div> <div>93%</div> <div>0%</div> </div>
1	B	343	<div> <div>2%</div> <div>92%</div> <div>6%</div> </div>
1	C	343	<div> <div>3%</div> <div>88%</div> <div>7%</div> </div>
1	D	343	<div> <div>12%</div> <div>89%</div> <div>8%</div> </div>

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 12387 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 2,3-dihydroxybenzoate decarboxylase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	334	Total	C	N	O	S	0	0	0
			2692	1737	460	486	9			
1	B	339	Total	C	N	O	S	0	0	0
			2739	1764	471	495	9			
1	C	328	Total	C	N	O	S	0	0	0
			2638	1701	452	476	9			
1	D	332	Total	C	N	O	S	0	0	0
			2677	1729	457	482	9			

There are 28 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	HIS	-	expression tag	UNP A0A420U2F4
A	338	HIS	-	expression tag	UNP A0A420U2F4
A	339	HIS	-	expression tag	UNP A0A420U2F4
A	340	HIS	-	expression tag	UNP A0A420U2F4
A	341	HIS	-	expression tag	UNP A0A420U2F4
A	342	HIS	-	expression tag	UNP A0A420U2F4
A	343	HIS	-	expression tag	UNP A0A420U2F4
B	1	HIS	-	expression tag	UNP A0A420U2F4
B	338	HIS	-	expression tag	UNP A0A420U2F4
B	339	HIS	-	expression tag	UNP A0A420U2F4
B	340	HIS	-	expression tag	UNP A0A420U2F4
B	341	HIS	-	expression tag	UNP A0A420U2F4
B	342	HIS	-	expression tag	UNP A0A420U2F4
B	343	HIS	-	expression tag	UNP A0A420U2F4
C	1	HIS	-	expression tag	UNP A0A420U2F4
C	338	HIS	-	expression tag	UNP A0A420U2F4
C	339	HIS	-	expression tag	UNP A0A420U2F4
C	340	HIS	-	expression tag	UNP A0A420U2F4
C	341	HIS	-	expression tag	UNP A0A420U2F4
C	342	HIS	-	expression tag	UNP A0A420U2F4
C	343	HIS	-	expression tag	UNP A0A420U2F4

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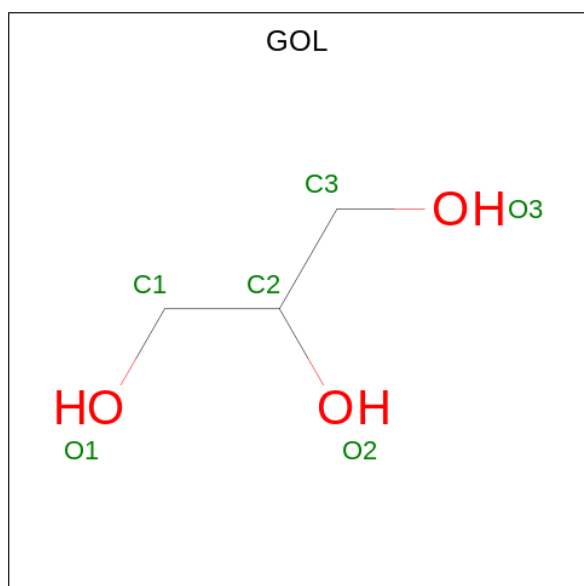
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Chain	Residue	Modelled	Actual	Comment	Reference
D	1	HIS	-	expression tag	UNP A0A420U2F4
D	338	HIS	-	expression tag	UNP A0A420U2F4
D	339	HIS	-	expression tag	UNP A0A420U2F4
D	340	HIS	-	expression tag	UNP A0A420U2F4
D	341	HIS	-	expression tag	UNP A0A420U2F4
D	342	HIS	-	expression tag	UNP A0A420U2F4
D	343	HIS	-	expression tag	UNP A0A420U2F4

- 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	1	Total	Zn	0	0
			1	1		
2	A	1	Total	Zn	0	0
			1	1		
2	D	1	Total	Zn	0	0
			1	1		
2	C	1	Total	Zn	0	0
			1	1		

- Molecule 3 is GLYCEROL (three-letter code: GOL) (formula: C₃H₈O₃).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	B	1	Total	C	O	0	0
			6	3	3		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	B	1	Total	C	O	0	0
			6	3	3		
3	C	1	Total	C	O	0	0
			6	3	3		

- Molecule 4 is water.

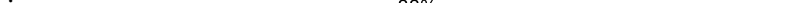
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	450	Total	O	0	0
			450	450		
4	B	477	Total	O	0	0
			477	477		
4	C	412	Total	O	0	0
			412	412		
4	D	280	Total	O	0	0
			280	280		

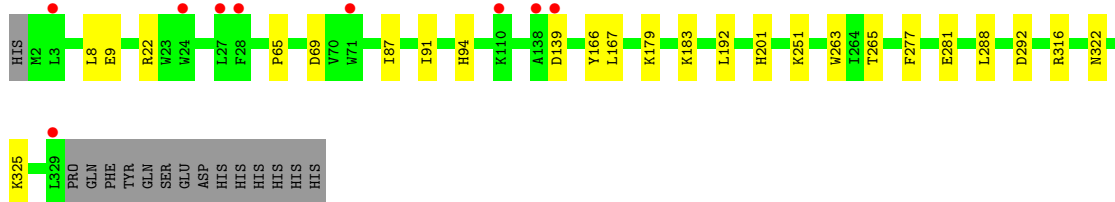
- Molecule 1: 2,3-dihydroxybenzoate decarboxylase

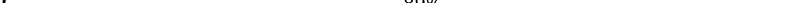


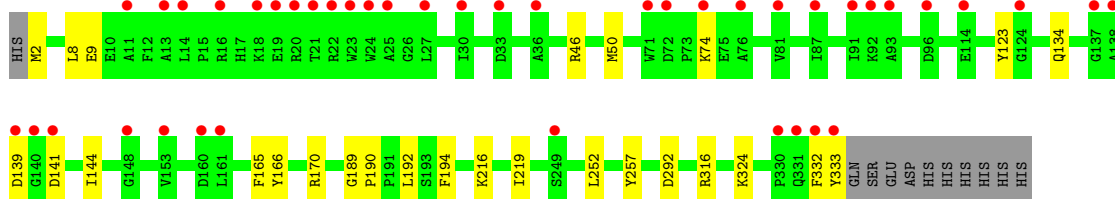
- Chain B:  2% 92% 6%



- Chain C:  3% 88% 7%



- Chain D:  12% 89% 8%



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	77.62Å 132.25Å 140.59Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	24.77 – 1.55 24.77 – 1.55	Depositor EDS
% Data completeness (in resolution range)	99.0 (24.77-1.55) 99.0 (24.77-1.55)	Depositor EDS
R_{merge}	0.04	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.73 (at 1.55Å)	Xtriage
Refinement program	PHENIX 1.17.1 _3660	Depositor
R, R_{free}	0.186 , 0.199 0.190 , 0.202	Depositor DCC
R_{free} test set	10338 reflections (4.99%)	wwPDB-VP
Wilson B-factor (Å ²)	20.3	Xtriage
Anisotropy	0.180	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 49.8	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	0.96	EDS
Total number of atoms	12387	wwPDB-VP
Average B, all atoms (Å ²)	26.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.23% 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: GOL, 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	0.36	0/2773	0.58	0/3767
1	B	0.35	0/2823	0.58	0/3835
1	C	0.35	0/2716	0.57	0/3689
1	D	0.33	0/2758	0.53	0/3747
All	All	0.35	0/11070	0.57	0/15038

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 [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	2692	0	2611	8	0
1	B	2739	0	2642	14	0
1	C	2638	0	2565	13	0
1	D	2677	0	2598	17	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	B	12	0	15	2	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	C	6	0	7	0	0
4	A	450	0	0	3	0
4	B	477	0	0	3	0
4	C	412	0	0	4	0
4	D	280	0	0	2	0
All	All	12387	0	10438	49	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 2.

The worst 5 of 49 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:92:LYS:HE3	4:A:816:HOH:O	1.98	0.62
1:B:267:SER:HB2	1:B:292:ASP:HB2	1.84	0.60
1:D:2:MET:CE	1:D:316:ARG:HA	2.36	0.54
1:B:52:GLU:HG2	1:B:340:HIS:NE2	2.24	0.53
1:D:74:LYS:H	1:D:74:LYS:HE2	1.73	0.53

There are no symmetry-related clashes.

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	332/343 (97%)	320 (96%)	11 (3%)	1 (0%)	41	19
1	B	337/343 (98%)	326 (97%)	10 (3%)	1 (0%)	41	19
1	C	326/343 (95%)	315 (97%)	10 (3%)	1 (0%)	41	19
1	D	330/343 (96%)	315 (96%)	14 (4%)	1 (0%)	41	19
All	All	1325/1372 (97%)	1276 (96%)	45 (3%)	4 (0%)	41	19

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	192	LEU
1	B	192	LEU
1	C	192	LEU
1	D	192	LEU

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	281/290 (97%)	278 (99%)	3 (1%)	73	53
1	B	286/290 (99%)	282 (99%)	4 (1%)	67	41
1	C	275/290 (95%)	271 (98%)	4 (2%)	65	37
1	D	279/290 (96%)	275 (99%)	4 (1%)	67	41
All	All	1121/1160 (97%)	1106 (99%)	15 (1%)	69	44

5 of 15 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	340	HIS
1	C	8	LEU
1	D	139	ASP
1	B	170	ARG
1	D	8	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	D	331	GLN

5.3.3 RNA ⓘ

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 7 ligands modelled in this entry, 4 are monoatomic - leaving 3 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	$\# Z > 2$
3	GOL	B	402	-	5,5,5	1.07	0	5,5,5	1.11	0
3	GOL	C	401	-	5,5,5	1.72	2 (40%)	5,5,5	0.68	0
3	GOL	B	401	-	5,5,5	1.38	1 (20%)	5,5,5	0.96	0

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
3	GOL	B	402	-	-	0/4/4/4	-
3	GOL	C	401	-	-	0/4/4/4	-
3	GOL	B	401	-	-	2/4/4/4	-

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	C	401	GOL	O2-C2	-2.71	1.35	1.43
3	B	401	GOL	O2-C2	-2.34	1.36	1.43
3	C	401	GOL	C1-C2	2.09	1.60	1.51

There are no bond angle outliers.

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	B	401	GOL	O1-C1-C2-C3
3	B	401	GOL	O1-C1-C2-O2

There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	402	GOL	1	0
3	B	401	GOL	1	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	334/343 (97%)	-0.20	5 (1%) 73 78	13, 19, 30, 54	0
1	B	339/343 (98%)	-0.06	8 (2%) 59 65	14, 22, 35, 47	0
1	C	328/343 (95%)	0.12	9 (2%) 54 62	14, 24, 36, 47	0
1	D	332/343 (96%)	0.74	42 (12%) 3 3	17, 31, 46, 63	0
All	All	1333/1372 (97%)	0.15	64 (4%) 30 35	13, 23, 40, 63	0

The worst 5 of 64 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	138	ALA	10.3
1	C	138	ALA	9.7
1	D	333	TYR	9.2
1	D	23	TRP	8.1
1	D	331	GLN	7.1

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](#)

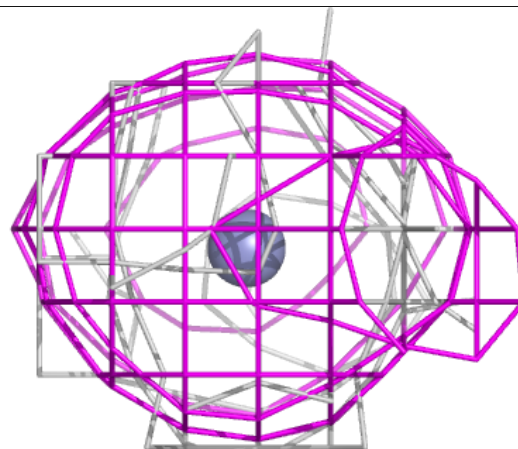
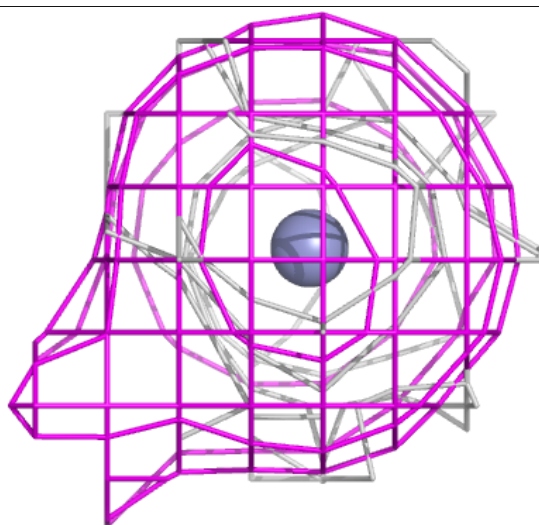
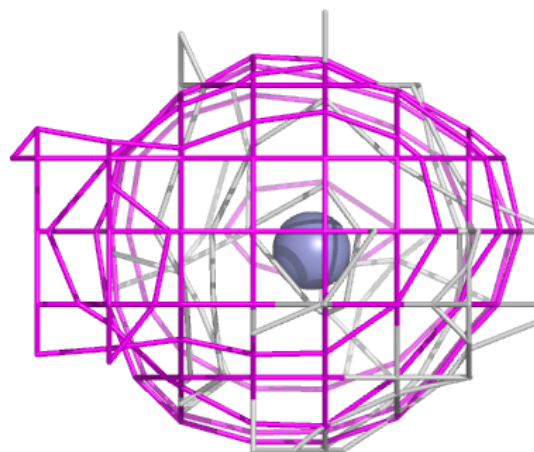
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled ‘Q< 0.9’ lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
3	GOL	B	401	6/6	0.87	0.12	26,29,29,33	0
2	ZN	A	401	1/1	0.90	0.52	30,30,30,30	0
3	GOL	B	402	6/6	0.90	0.12	24,24,30,32	0
2	ZN	B	403	1/1	0.94	0.44	30,30,30,30	0
2	ZN	C	402	1/1	0.95	0.42	30,30,30,30	0
3	GOL	C	401	6/6	0.96	0.09	16,19,22,24	0
2	ZN	D	401	1/1	0.98	0.57	30,30,30,30	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.

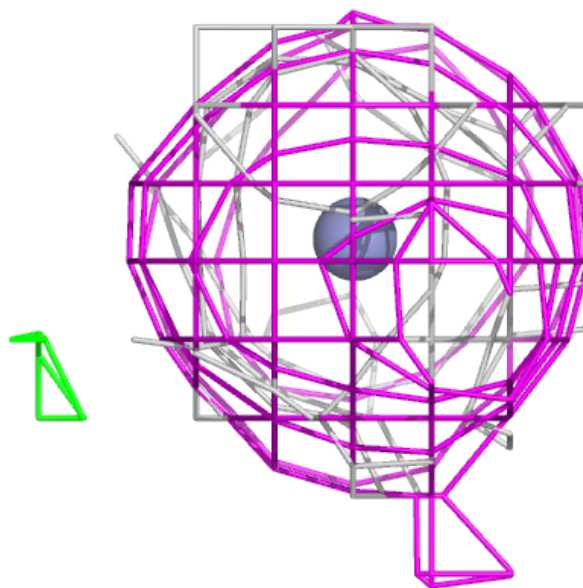
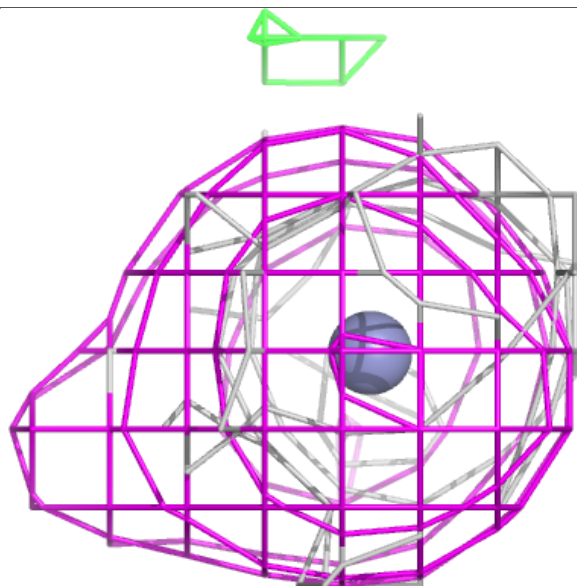
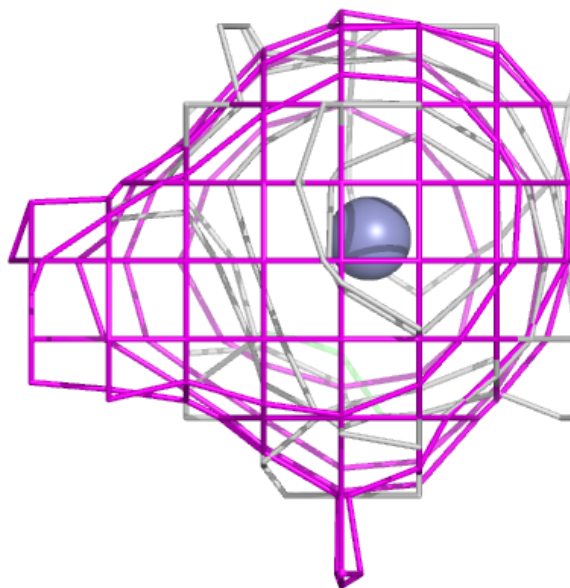
Electron density around ZN A 401:

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



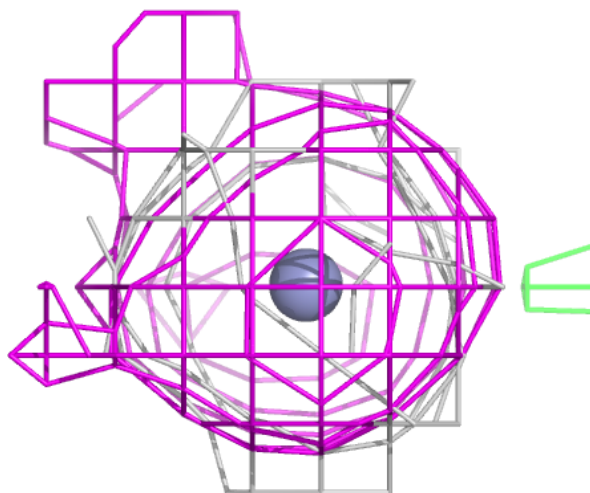
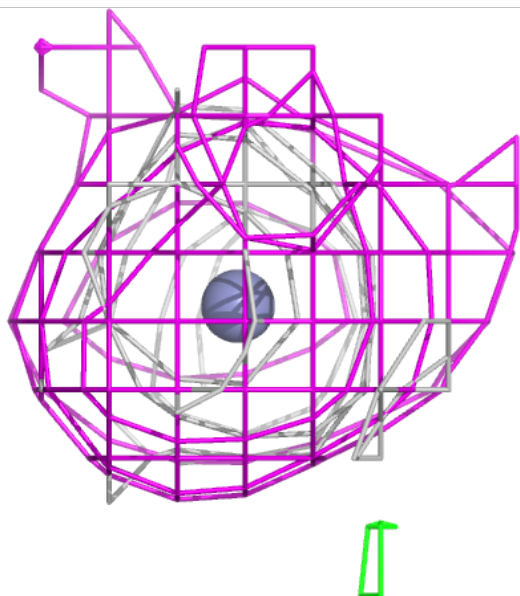
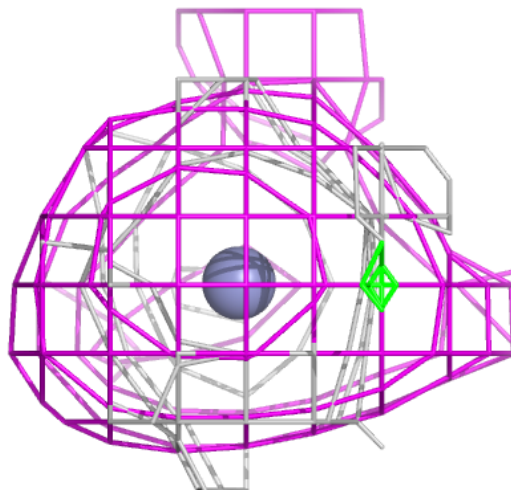
Electron density around ZN B 403:

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



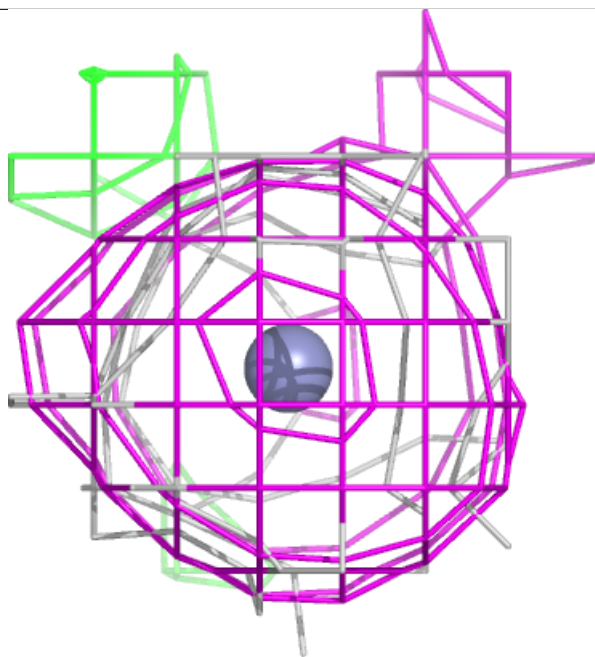
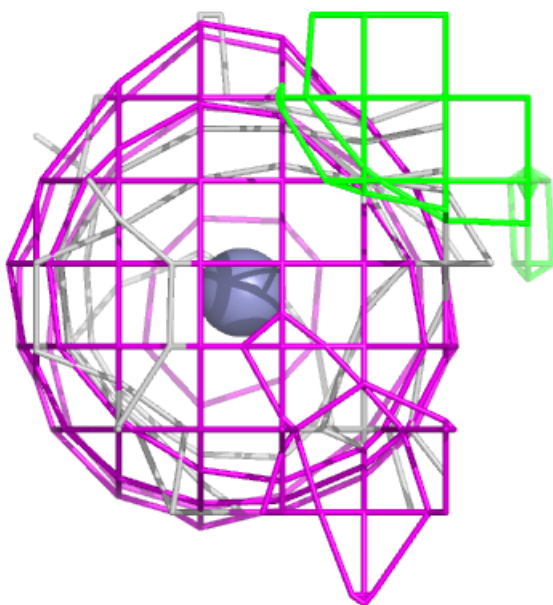
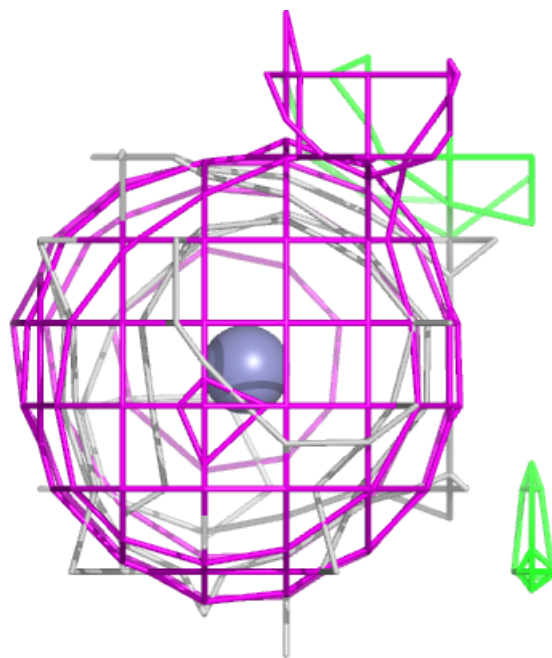
Electron density around ZN C 402:

$2mF_o-DF_c$ (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 [i](#)

There are no such residues in this entry.