



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 24, 2025 – 12:15 pm GMT

PDB ID : 9H9Z  
Title : Crystal structure of Cu(II)-bound LmrR\_V15Bpy  
Authors : Thunnissen, A.M.W.H.; Jiang, R.; Casilli, F.; Aalbers, F.; Roelfes, G.  
Deposited on : 2024-11-01  
Resolution : 1.90 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](mailto: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>.

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

MolProbity	:	4.02b-467
Mogul	:	1.8.4, CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	3.0
Percentile statistics	:	20231227.v01 (using entries in the PDB archive December 27th 2023)
CCP4	:	9.0.003 (Gargrove)
Density-Fitness	:	1.0.11
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.41

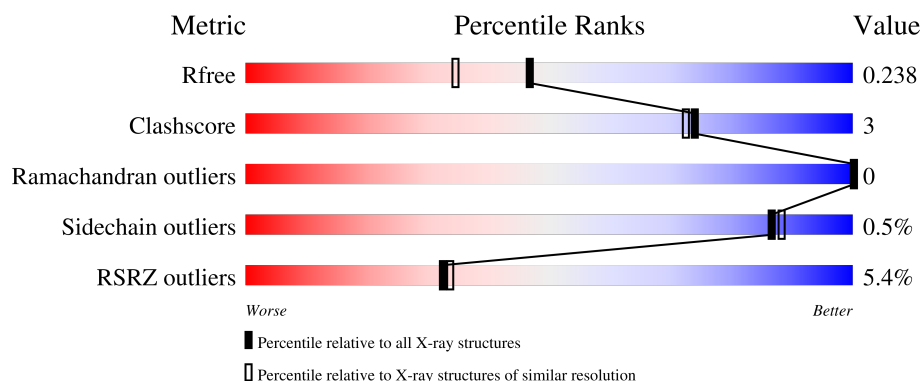
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 1.90 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))
$R_{free}$	164625	7293 (1.90-1.90)
Clashscore	180529	8090 (1.90-1.90)
Ramachandran outliers	177936	8022 (1.90-1.90)
Sidechain outliers	177891	8022 (1.90-1.90)
RSRZ outliers	164620	7292 (1.90-1.90)

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  $\geq 3$ , 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	131	<div> <div>5%</div> <div>78%</div> <div>8%</div> <div>14%</div> </div>
1	B	131	<div> <div>5%</div> <div>79%</div> <div>6%</div> <div>15%</div> </div>

## 2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 3784 atoms, of which 1841 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 Transcriptional regulator, PadR-like family.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	113	Total	C	H	N	O	S	0	1	0
			1866	592	931	164	176	3			
1	B	112	Total	C	H	N	O	S	0	0	0
			1829	582	908	161	175	3			

There are 34 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	GLY	-	expression tag	UNP A2RI36
A	15	BP5	VAL	engineered mutation	UNP A2RI36
A	117	SER	-	expression tag	UNP A2RI36
A	118	ARG	-	expression tag	UNP A2RI36
A	119	GLY	-	expression tag	UNP A2RI36
A	120	GLY	-	expression tag	UNP A2RI36
A	121	SER	-	expression tag	UNP A2RI36
A	122	GLY	-	expression tag	UNP A2RI36
A	123	GLY	-	expression tag	UNP A2RI36
A	124	TRP	-	expression tag	UNP A2RI36
A	125	SER	-	expression tag	UNP A2RI36
A	126	HIS	-	expression tag	UNP A2RI36
A	127	PRO	-	expression tag	UNP A2RI36
A	128	GLN	-	expression tag	UNP A2RI36
A	129	PHE	-	expression tag	UNP A2RI36
A	130	GLU	-	expression tag	UNP A2RI36
A	131	LYS	-	expression tag	UNP A2RI36
B	1	GLY	-	expression tag	UNP A2RI36
B	15	BP5	VAL	engineered mutation	UNP A2RI36
B	117	SER	-	expression tag	UNP A2RI36
B	118	ARG	-	expression tag	UNP A2RI36
B	119	GLY	-	expression tag	UNP A2RI36
B	120	GLY	-	expression tag	UNP A2RI36
B	121	SER	-	expression tag	UNP A2RI36
B	122	GLY	-	expression tag	UNP A2RI36

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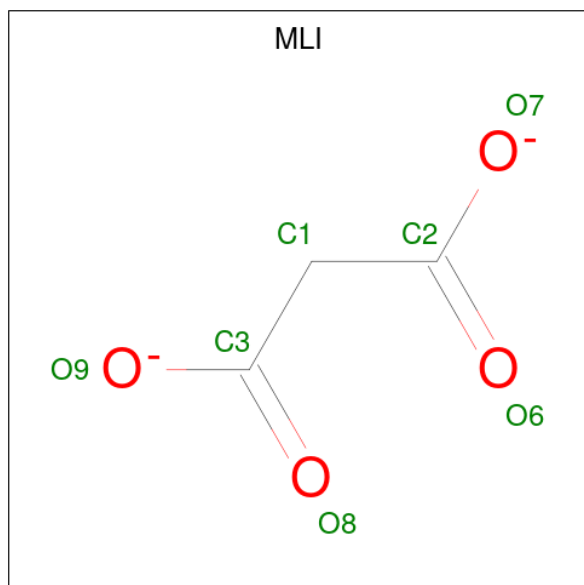
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Chain	Residue	Modelled	Actual	Comment	Reference
B	123	GLY	-	expression tag	UNP A2RI36
B	124	TRP	-	expression tag	UNP A2RI36
B	125	SER	-	expression tag	UNP A2RI36
B	126	HIS	-	expression tag	UNP A2RI36
B	127	PRO	-	expression tag	UNP A2RI36
B	128	GLN	-	expression tag	UNP A2RI36
B	129	PHE	-	expression tag	UNP A2RI36
B	130	GLU	-	expression tag	UNP A2RI36
B	131	LYS	-	expression tag	UNP A2RI36

- Molecule 2 is COPPER (II) ION (three-letter code: CU) (formula: Cu) (labeled as "Ligand of Interest" by depositor).

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

- Molecule 3 is MALONATE ION (three-letter code: MLI) (formula: C<sub>3</sub>H<sub>2</sub>O<sub>4</sub>).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	B	1	Total C H O 9 3 2 4	0	0

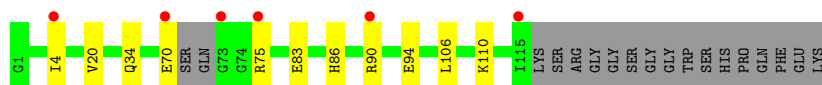
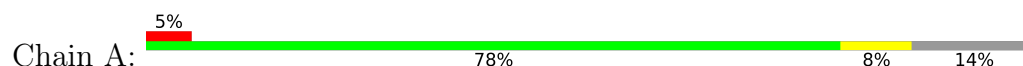
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	41	Total 41	O 41	0	0
4	B	37	Total 37	O 37	0	0

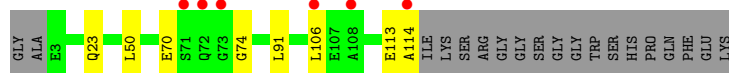
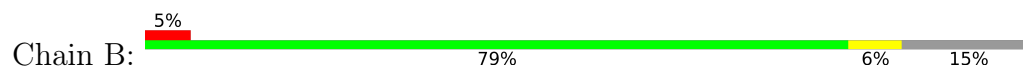
### 3 Residue-property plots [i](#)

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: Transcriptional regulator, PadR-like family



- Molecule 1: Transcriptional regulator, PadR-like family



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	35.50Å 53.36Å 147.78Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	34.51 – 1.90 34.51 – 1.90	Depositor EDS
% Data completeness (in resolution range)	98.1 (34.51-1.90) 98.0 (34.51-1.90)	Depositor EDS
$R_{merge}$	0.07	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.49 (at 1.89Å)	Xtriage
Refinement program	PHENIX 1.21.1_5286	Depositor
R, $R_{free}$	0.202 , 0.239 0.201 , 0.238	Depositor DCC
$R_{free}$ test set	1131 reflections (4.92%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	38.4	Xtriage
Anisotropy	0.259	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.40 , 34.8	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.96	EDS
Total number of atoms	3784	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	53.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.53% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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: CU, MLI, BP5

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	0.64	1/932 (0.1%)	0.72	0/1245
1	B	0.59	0/916	0.73	0/1226
All	All	0.62	1/1848 (0.1%)	0.73	0/2471

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	83	GLU	CB-CG	5.12	1.61	1.52

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	935	931	934	9	0
1	B	921	908	909	8	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	B	7	2	2	0	0
4	A	41	0	0	0	0
4	B	37	0	0	0	0
All	All	1943	1841	1845	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 3.

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:A:90[B]:ARG:NH1	1:A:94:GLU:OE2	2.41	0.53
1:B:70:GLU:HA	1:B:74:GLY:O	2.11	0.51
1:A:34:GLN:HB3	1:B:106:LEU:HD21	1.93	0.49
1:A:4:ILE:HD11	1:B:91:LEU:HB3	1.94	0.49
1:A:106:LEU:HB3	1:B:23:GLN:HE22	1.79	0.48
1:A:70:GLU:HG3	1:A:75:ARG:HH21	1.81	0.46
1:A:20:VAL:HG22	1:B:106:LEU:CD1	2.46	0.45
1:A:86:HIS:HB3	1:A:90[A]:ARG:NH2	2.32	0.44
1:A:20:VAL:HG22	1:B:106:LEU:HD13	2.01	0.42
1:A:110:LYS:HD2	1:B:23:GLN:HB3	2.02	0.42
1:B:113:GLU:O	1:B:114:ALA:O	2.39	0.41

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	109/131 (83%)	107 (98%)	2 (2%)	0	100	100
1	B	109/131 (83%)	107 (98%)	2 (2%)	0	100	100
All	All	218/262 (83%)	214 (98%)	4 (2%)	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	97/110 (88%)	97 (100%)	0	100	100
1	B	96/110 (87%)	95 (99%)	1 (1%)	73	74
All	All	193/220 (88%)	192 (100%)	1 (0%)	86	88

All (1) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	50	LEU

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

### 5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

2 non-standard protein/DNA/RNA residues are modelled in this entry.

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	$\# Z  > 2$
1	BP5	A	15	2,1	17,18,19	0.43	0	20,23,25	0.71	0
1	BP5	B	15	2,1	17,18,19	0.47	0	20,23,25	0.47	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
1	BP5	A	15	2,1	-	4/9/10/12	0/2/2/2
1	BP5	B	15	2,1	-	3/9/10/12	0/2/2/2

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (7) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	A	15	BP5	N1-C3-C6-N2
1	B	15	BP5	C2-C3-C6-N2
1	A	15	BP5	N1-C3-C6-C7
1	A	15	BP5	C2-C3-C6-C7
1	B	15	BP5	C2-C3-C6-C7
1	B	15	BP5	N1-C3-C6-N2
1	A	15	BP5	C2-C3-C6-N2

There are no ring outliers.

No monomer is involved in short contacts.

## 5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

## 5.6 Ligand geometry [i](#)

Of 3 ligands modelled in this entry, 2 are monoatomic - leaving 1 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	MLI	B	201	2	6,6,6	1.97	1 (16%)	7,7,7	0.48	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	MLI	B	201	2	-	2/4/4/4	-

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	201	MLI	C1-C3	3.47	1.56	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	201	MLI	C3-C1-C2-O6
3	B	201	MLI	C3-C1-C2-O7

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<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	112/131 (85%)	0.27	6 (5%) 32 34	31, 50, 92, 105	1 (0%)
1	B	111/131 (84%)	0.18	6 (5%) 32 34	34, 47, 87, 127	0
All	All	223/262 (85%)	0.23	12 (5%) 32 34	31, 48, 91, 127	1 (0%)

All (12) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	114	ALA	4.9
1	A	115	ILE	4.5
1	B	72	GLN	3.6
1	A	73	GLY	3.6
1	B	108	ALA	3.0
1	B	106	LEU	2.9
1	B	71	SER	2.5
1	A	70	GLU	2.5
1	B	73	GLY	2.4
1	A	75	ARG	2.3
1	A	4	ILE	2.3
1	A	90[A]	ARG	2.0

### 6.2 Non-standard residues in protein, DNA, RNA chains [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<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	BP5	A	15	17/18	0.94	0.10	31,45,77,93	0
1	BP5	B	15	17/18	0.94	0.11	29,42,58,65	0

### 6.3 Carbohydrates [i](#)

There are no monosaccharides in this entry.

### 6.4 Ligands [i](#)

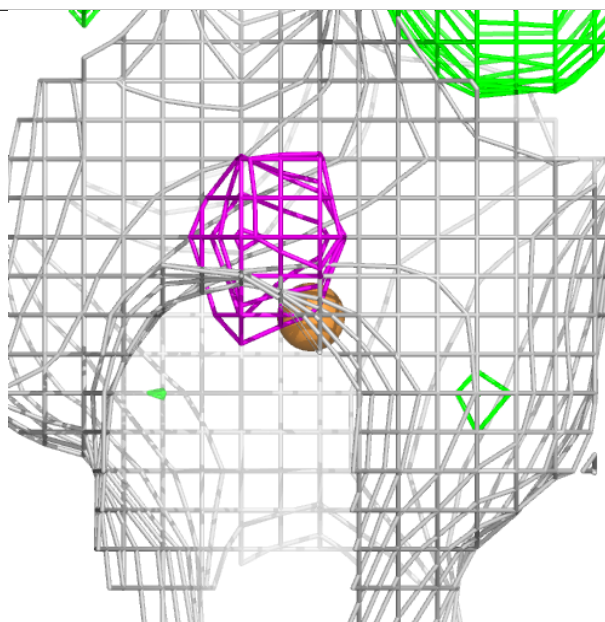
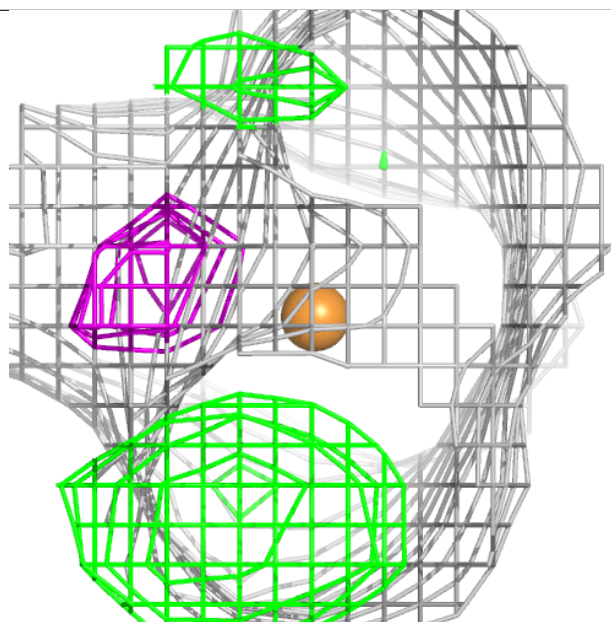
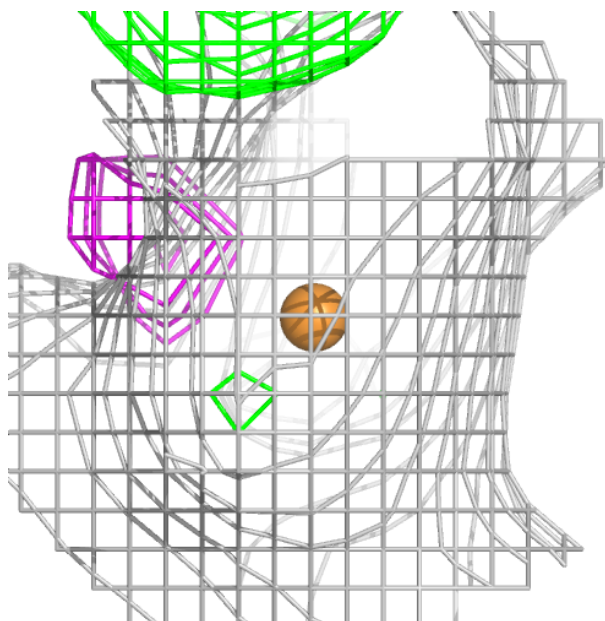
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 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
3	MLI	B	201	7/7	0.87	0.12	44,53,68,69	0
2	CU	A	201	1/1	0.93	0.08	81,81,81,81	0
2	CU	B	202	1/1	0.99	0.09	56,56,56,56	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 CU A 201:**

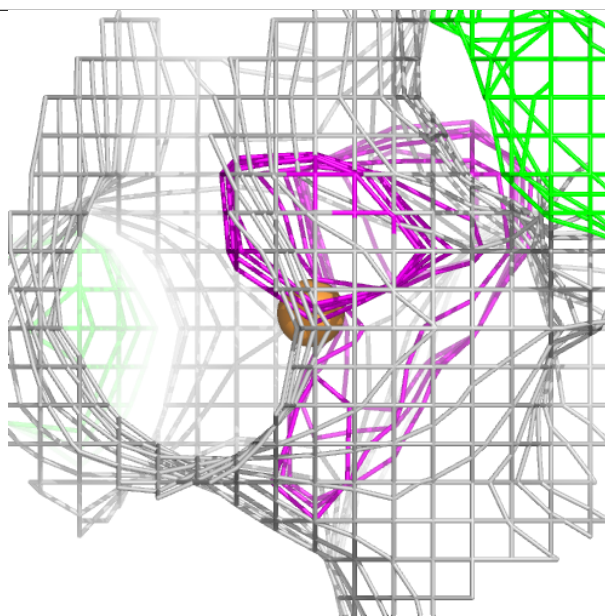
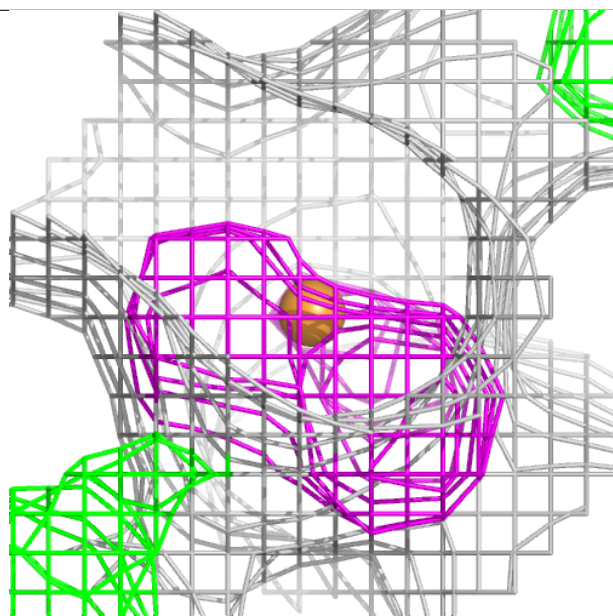
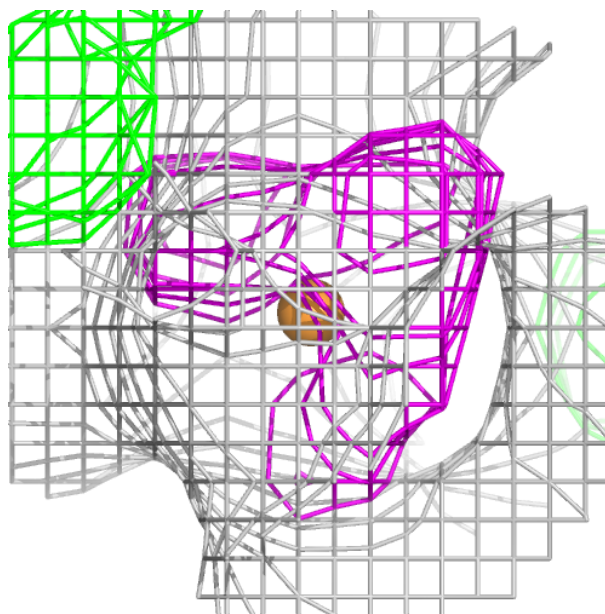
$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)





**Electron density around CU B 202:**

$2mF_o-DF_c$  (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.