



# Full wwPDB X-ray Structure Validation Report ⓘ

Aug 29, 2020 – 04:23 PM BST

PDB ID : 6GYI  
Title : Azurin fom Pseudomonas aeruginosa treated with hydrosulfide  
Authors : Palm, G.J.; Kohlhouse, D.  
Deposited on : 2018-06-29  
Resolution : 1.60 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.

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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  
Xtriage (Phenix) : 1.13  
EDS : 2.13  
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) : 2.13

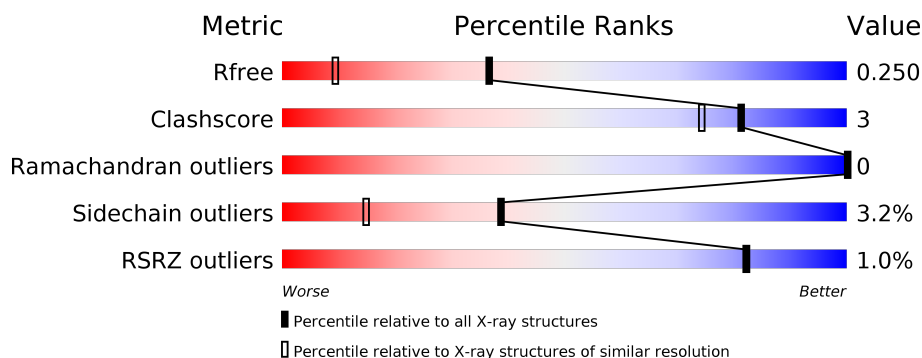
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 1.60 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	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
$R_{free}$	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 on 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	128	 85% 13% ...
1	B	128	 86% 12% ..
1	C	128	 84% 16% .
1	D	128	 72% 27% ..

## 2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 4331 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 Azurin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	127	Total	C	N	O	S	0	2	0
			971	609	162	190	10			
1	B	127	Total	C	N	O	S	0	3	0
			975	610	163	193	9			
1	C	128	Total	C	N	O	S	0	2	0
			974	609	163	193	9			
1	D	127	Total	C	N	O	S	0	2	0
			976	610	163	194	9			

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

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

- Molecule 3 is CALCIUM ION (three-letter code: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	1	Total	Ca	0	0
			1	1		
3	A	1	Total	Ca	0	0
			1	1		
3	D	1	Total	Ca	0	0
			1	1		
3	C	1	Total	Ca	0	0
			1	1		


- Molecule 4 is water.

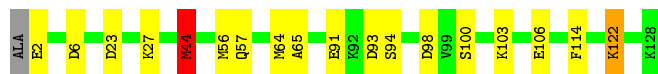
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	110	Total 110	O 110	0	0
4	B	97	Total 97	O 97	0	0
4	C	106	Total 106	O 106	0	0
4	D	114	Total 114	O 114	0	0

### 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.

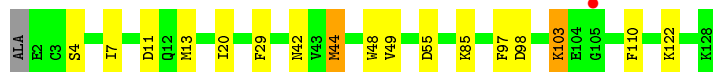
- Molecule 1: Azurin

Chain A: 




- Molecule 1: Azurin

Chain B: 



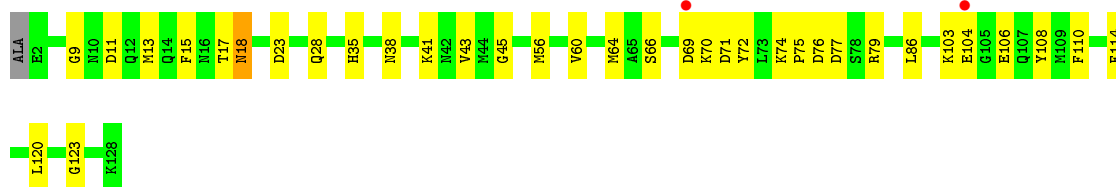
- Molecule 1: Azurin

Chain C: 



- Molecule 1: Azurin

Chain D: 



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	50.74Å 59.78Å 79.73Å 90.00° 90.38° 90.00°	Depositor
Resolution (Å)	47.83 – 1.60 47.83 – 1.60	Depositor EDS
% Data completeness (in resolution range)	99.2 (47.83-1.60) 99.2 (47.83-1.60)	Depositor EDS
$R_{merge}$	0.03	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.16 (at 1.60Å)	Xtriage
Refinement program	REFMAC 5.8.0189	Depositor
R, $R_{free}$	0.192 , 0.243 0.198 , 0.250	Depositor DCC
$R_{free}$ test set	2089 reflections (3.34%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	20.0	Xtriage
Anisotropy	0.816	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.33 , 39.0	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.080 for h,-k,-l	Xtriage
$F_o, F_c$ correlation	0.96	EDS
Total number of atoms	4331	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	30.0	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 27.52 % 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 2.1687e-03. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

<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: CA, CU

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	1.49	5/994 (0.5%)	1.50	15/1341 (1.1%)
1	B	1.42	2/1001 (0.2%)	1.37	8/1350 (0.6%)
1	C	1.36	2/997 (0.2%)	1.37	11/1345 (0.8%)
1	D	1.70	11/999 (1.1%)	1.58	15/1348 (1.1%)
All	All	1.50	20/3991 (0.5%)	1.46	49/5384 (0.9%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	D	0	1

All (20) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	108	TYR	CE2-CZ	-8.71	1.27	1.38
1	A	106	GLU	CD-OE1	8.47	1.34	1.25
1	B	97	PHE	CG-CD2	-7.59	1.27	1.38
1	A	100	SER	CB-OG	7.45	1.51	1.42
1	D	108	TYR	CZ-OH	7.11	1.50	1.37
1	B	4	SER	CB-OG	-6.51	1.33	1.42
1	A	94	SER	CB-OG	-6.45	1.33	1.42
1	A	65	ALA	C-O	-5.91	1.12	1.23
1	D	15	PHE	CG-CD1	-5.90	1.29	1.38
1	D	43	VAL	C-O	5.83	1.34	1.23
1	D	106	GLU	CD-OE1	5.70	1.31	1.25
1	A	2	GLU	C-O	5.70	1.34	1.23
1	D	69	ASP	CB-CG	5.69	1.63	1.51
1	D	45	GLY	CA-C	-5.68	1.42	1.51

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	110	PHE	CE1-CZ	5.58	1.48	1.37
1	D	123	GLY	C-O	-5.41	1.15	1.23
1	D	60	VAL	C-O	5.38	1.33	1.23
1	D	41	LYS	CA-CB	-5.27	1.42	1.53
1	C	123	GLY	C-O	-5.18	1.15	1.23
1	C	89	SER	CB-OG	-5.05	1.35	1.42

All (49) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	56	MET	CG-SD-CE	-10.12	84.02	100.20
1	C	6	ASP	CB-CG-OD1	8.93	126.34	118.30
1	A	44[A]	MET	CA-CB-CG	8.73	128.15	113.30
1	A	44[B]	MET	CA-CB-CG	8.73	128.15	113.30
1	C	6	ASP	CB-CG-OD2	-8.52	110.63	118.30
1	A	64	MET	CG-SD-CE	8.50	113.79	100.20
1	A	93	ASP	CB-CG-OD2	8.28	125.75	118.30
1	B	98	ASP	CB-CG-OD1	7.96	125.46	118.30
1	D	23	ASP	CB-CG-OD1	7.84	125.36	118.30
1	D	114	PHE	CB-CG-CD1	-7.76	115.37	120.80
1	D	71	ASP	CB-CG-OD2	-7.45	111.59	118.30
1	A	6	ASP	CB-CG-OD1	7.34	124.91	118.30
1	D	13	MET	CG-SD-CE	7.26	111.82	100.20
1	D	11	ASP	CB-CG-OD1	6.67	124.30	118.30
1	C	98	ASP	CB-CG-OD2	-6.64	112.33	118.30
1	D	108	TYR	CB-CG-CD2	-6.44	117.14	121.00
1	D	38	ASN	N-CA-CB	-6.43	99.03	110.60
1	B	11	ASP	CB-CG-OD1	6.33	124.00	118.30
1	D	64	MET	CG-SD-CE	6.29	110.26	100.20
1	A	122	LYS	CD-CE-NZ	-6.27	97.27	111.70
1	C	79	ARG	NE-CZ-NH2	6.21	123.41	120.30
1	A	6	ASP	CB-CG-OD2	-6.08	112.83	118.30
1	D	77	ASP	CB-CG-OD1	6.05	123.75	118.30
1	B	98	ASP	CB-CG-OD2	-6.04	112.86	118.30
1	A	56	MET	CG-SD-CE	-5.99	90.62	100.20
1	D	79	ARG	NE-CZ-NH1	-5.89	117.36	120.30
1	C	11	ASP	CB-CG-OD1	5.83	123.55	118.30
1	A	23	ASP	CB-CG-OD1	5.83	123.54	118.30
1	B	55	ASP	CB-CG-OD2	-5.76	113.12	118.30
1	B	122	LYS	CD-CE-NZ	-5.71	98.57	111.70
1	A	98	ASP	CB-CG-OD1	5.70	123.43	118.30
1	D	23	ASP	CB-CG-OD2	-5.69	113.18	118.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	76	ASP	CB-CG-OD2	5.55	123.30	118.30
1	C	98	ASP	CB-CG-OD1	5.52	123.26	118.30
1	A	23	ASP	CB-CG-OD2	-5.43	113.41	118.30
1	A	91	GLU	OE1-CD-OE2	-5.38	116.84	123.30
1	B	44	MET	CA-CB-CG	5.38	122.45	113.30
1	A	44[A]	MET	CB-CA-C	5.37	121.14	110.40
1	A	44[B]	MET	CB-CA-C	5.37	121.14	110.40
1	C	23	ASP	CB-CG-OD1	5.37	123.13	118.30
1	D	120	LEU	CB-CG-CD1	-5.27	102.04	111.00
1	C	108	TYR	CB-CG-CD1	5.25	124.15	121.00
1	C	11	ASP	CB-CG-OD2	-5.17	113.65	118.30
1	B	13	MET	CG-SD-CE	-5.08	92.08	100.20
1	C	115	PRO	CA-C-N	-5.07	106.07	116.20
1	B	29	PHE	CB-CG-CD2	-5.03	117.28	120.80
1	D	86	LEU	CB-CG-CD2	-5.03	102.45	111.00
1	C	13	MET	CG-SD-CE	-5.02	92.17	100.20
1	A	114	PHE	CB-CG-CD2	-5.01	117.29	120.80

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	D	72	TYR	Mainchain

## 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	971	0	951	4	0
1	B	975	0	953	6	0
1	C	974	0	953	5	0
1	D	976	0	952	6	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	A	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	1	0	0	0	0
3	C	1	0	0	0	0
3	D	1	0	0	0	0
4	A	110	0	0	2	0
4	B	97	0	0	0	0
4	C	106	0	0	0	0
4	D	114	0	0	1	0
All	All	4331	0	3809	20	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 (20) 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:7[A]:ILE:HD11	1:B:20:ILE:HD11	1.74	0.70
1:C:7[B]:ILE:HD11	1:C:20:ILE:HD11	1.74	0.69
1:D:17:THR:C	1:D:18:ASN:HD22	2.07	0.58
1:A:27:LYS:HD2	4:A:319:HOH:O	2.07	0.55
1:D:18:ASN:N	1:D:18:ASN:HD22	2.05	0.54
1:C:42:ASN:N	1:C:42:ASN:OD1	2.45	0.49
1:B:103:LYS:N	1:B:103:LYS:HD3	2.27	0.48
1:A:57:GLN:HB3	1:D:28:GLN:HE21	1.81	0.45
1:B:44:MET:HG3	1:B:44:MET:O	2.16	0.44
1:D:9:GLY:O	1:D:35:HIS:HA	2.18	0.43
1:B:49:VAL:O	1:B:110:PHE:HA	2.19	0.43
1:B:7[B]:ILE:HD13	1:B:7[B]:ILE:HG21	1.53	0.43
1:B:7[B]:ILE:CD1	1:B:48:TRP:CD1	3.02	0.42
1:D:74[B]:LYS:HD3	1:D:75:PRO:O	2.20	0.42
1:A:122:LYS:NZ	4:A:306:HOH:O	2.53	0.41
1:C:31:VAL:HG21	1:C:48:TRP:CH2	2.55	0.41
1:C:23:ASP:OD1	1:C:25:SER:OG	2.28	0.41
1:D:74[B]:LYS:NZ	4:D:303:HOH:O	2.48	0.41
1:C:20:ILE:HD12	1:C:110:PHE:CD1	2.56	0.40

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	127/128 (99%)	125 (98%)	2 (2%)	0	100	100
1	B	128/128 (100%)	125 (98%)	3 (2%)	0	100	100
1	C	128/128 (100%)	127 (99%)	1 (1%)	0	100	100
1	D	127/128 (99%)	124 (98%)	3 (2%)	0	100	100
All	All	510/512 (100%)	501 (98%)	9 (2%)	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	110/110 (100%)	107 (97%)	3 (3%)	44	20
1	B	111/110 (101%)	108 (97%)	3 (3%)	44	20
1	C	110/110 (100%)	106 (96%)	4 (4%)	35	12
1	D	111/110 (101%)	106 (96%)	5 (4%)	27	8
All	All	442/440 (100%)	427 (97%)	15 (3%)	39	13

All (15) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	44[A]	MET
1	A	44[B]	MET

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Mol	Chain	Res	Type
1	A	103	LYS
1	B	42	ASN
1	B	85	LYS
1	B	103	LYS
1	C	38	ASN
1	C	69	ASP
1	C	85	LYS
1	C	92	LYS
1	D	18	ASN
1	D	66	SER
1	D	70	LYS
1	D	103	LYS
1	D	104	GLU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	8	GLN
1	C	8	GLN
1	D	18	ASN
1	D	28	GLN

### 5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

## 5.5 Carbohydrates ⓘ

There are no monosaccharides in this entry.

## 5.6 Ligand geometry ⓘ

Of 8 ligands modelled in this entry, 8 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

There are no such residues in this entry.

## 5.8 Polymer linkage issues

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	127/128 (99%)	-0.19	0 <b>100</b> <b>100</b>	16, 26, 48, 59	0
1	B	127/128 (99%)	-0.23	1 (0%) <b>86</b> <b>86</b>	16, 27, 46, 58	0
1	C	128/128 (100%)	-0.06	2 (1%) <b>72</b> <b>71</b>	19, 34, 54, 68	0
1	D	127/128 (99%)	-0.06	2 (1%) <b>72</b> <b>71</b>	17, 27, 40, 62	0
All	All	509/512 (99%)	-0.14	5 (0%) <b>82</b> <b>82</b>	16, 28, 50, 68	0

All (5) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	105	GLY	4.8
1	D	104	GLU	2.5
1	C	38	ASN	2.4
1	C	105	GLY	2.1
1	D	69	ASP	2.1

### 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,

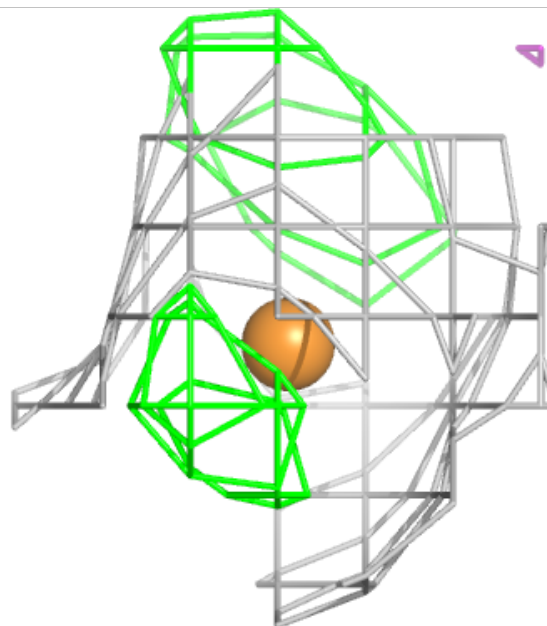
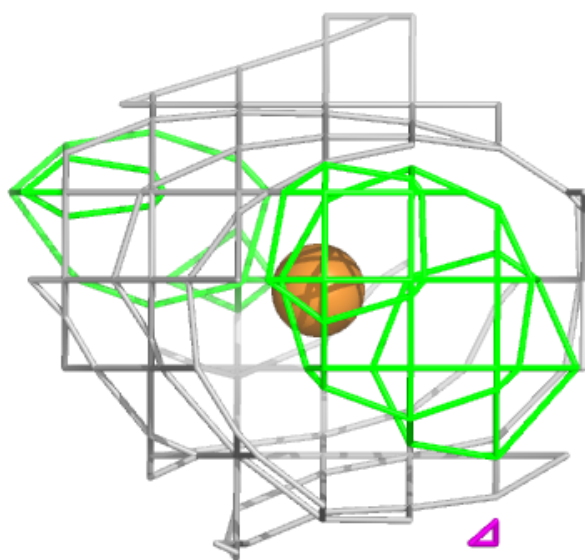
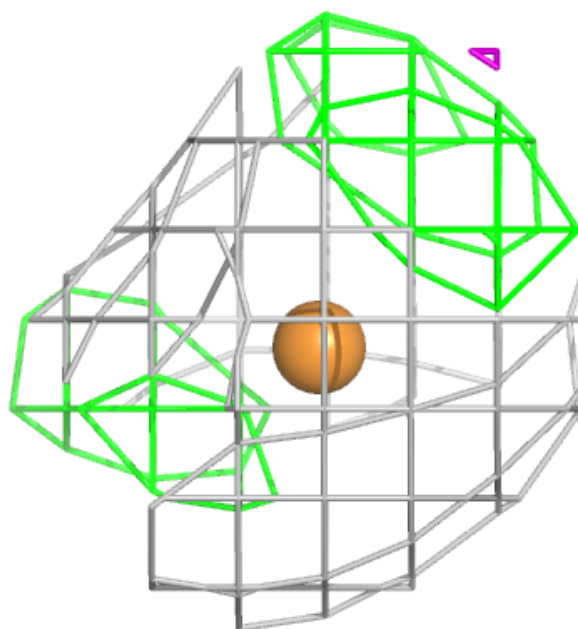
median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å <sup>2</sup> )	Q<0.9
3	CA	B	202	1/1	0.86	0.13	58,58,58,58	0
3	CA	C	202	1/1	0.90	0.13	61,61,61,61	0
2	CU	C	201	1/1	0.98	0.07	35,35,35,35	0
3	CA	D	202	1/1	0.99	0.08	30,30,30,30	0
2	CU	D	201	1/1	0.99	0.07	18,18,18,18	0
3	CA	A	202	1/1	0.99	0.06	30,30,30,30	0
2	CU	A	201	1/1	1.00	0.08	17,17,17,17	0
2	CU	B	201	1/1	1.00	0.07	22,22,22,22	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.

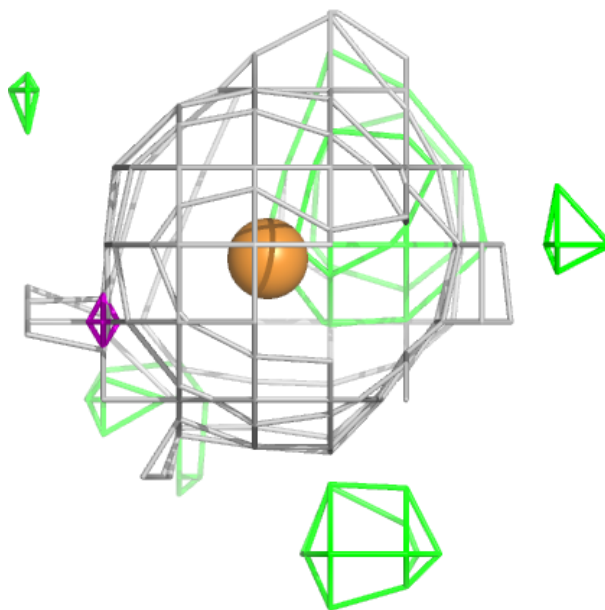
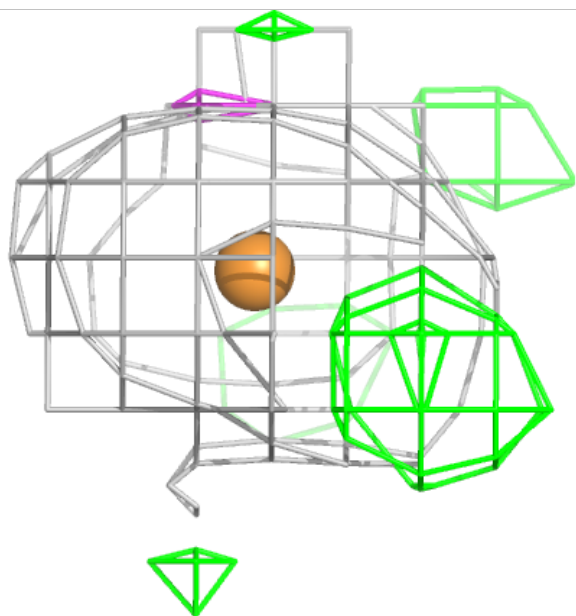
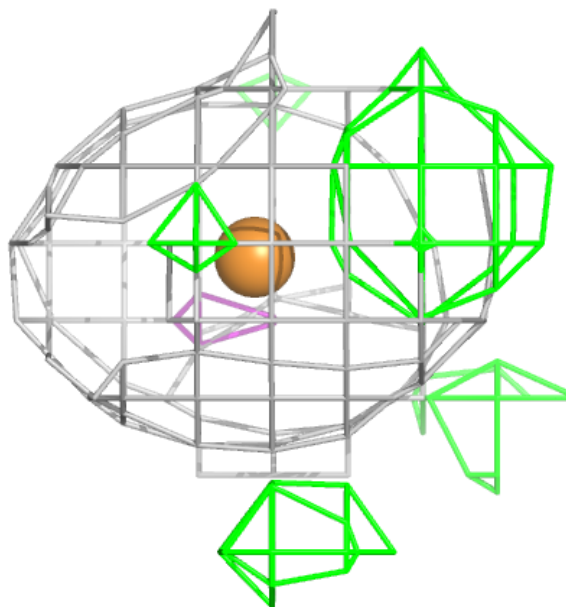
**Electron density around CU C 201:**

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



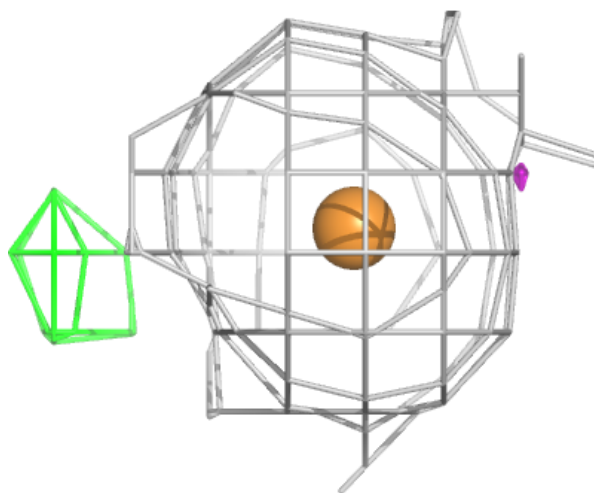
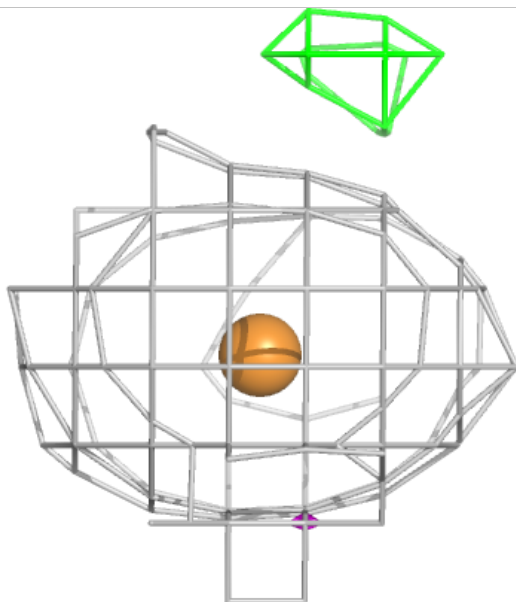
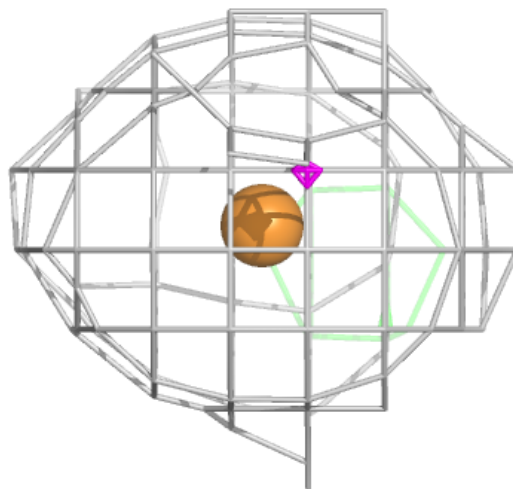
**Electron density around CU D 201:**

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



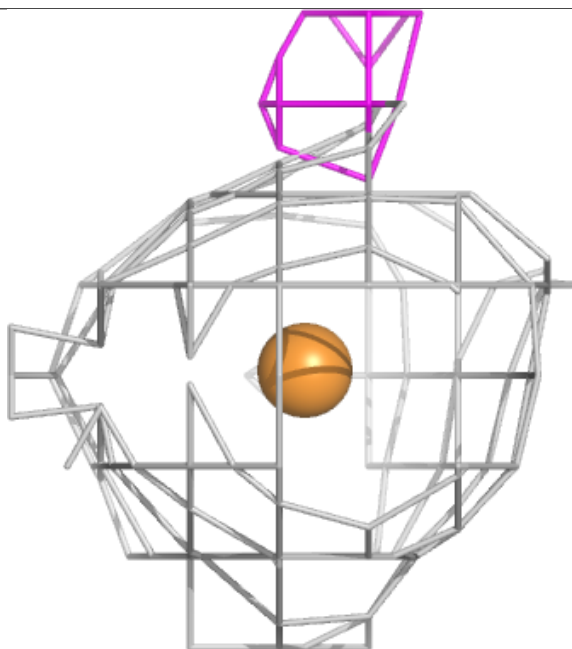
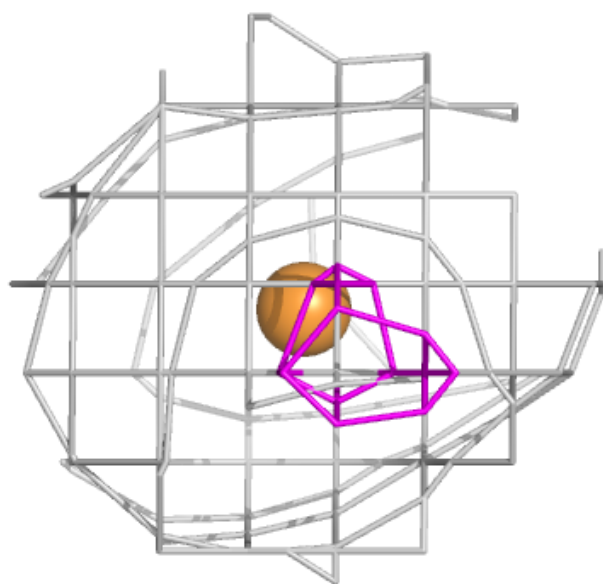
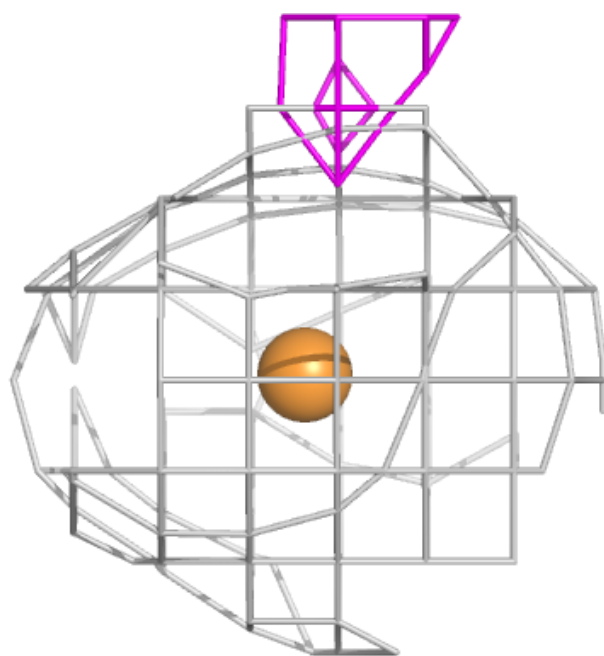
**Electron density around CU A 201:**

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



**Electron density around CU B 201:**

$2mF_o-DF_c$  (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.