



# Full wwPDB X-ray Structure Validation Report ⓘ

Oct 11, 2023 – 09:04 AM EDT

PDB ID : 7L5C  
Title : Structure of copper bound MEMO1  
Authors : Boniecki, M.T.; Uhlemann, E.E.; Dmitriev, O.Y.  
Deposited on : 2020-12-21  
Resolution : 2.55 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 : 1.8.5 (274361), CSD as541be (2020)  
Xtriage (Phenix) : 1.13  
EDS : 2.35.1  
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) : 2.35.1

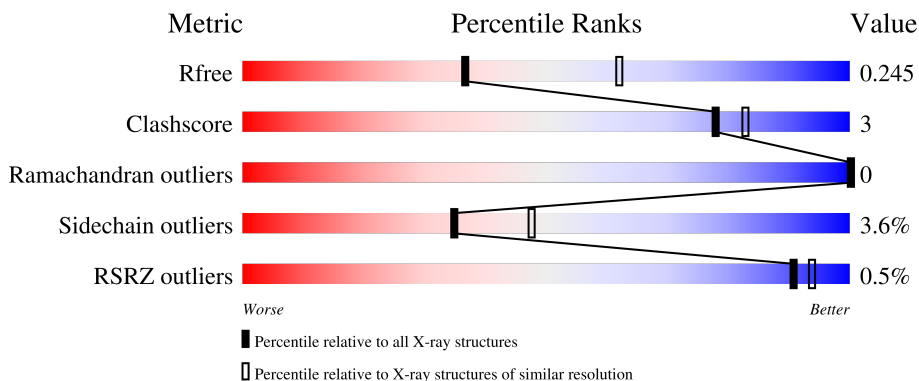
# 1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.55 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	1284 (2.56-2.52)
Clashscore	141614	1332 (2.56-2.52)
Ramachandran outliers	138981	1315 (2.56-2.52)
Sidechain outliers	138945	1315 (2.56-2.52)
RSRZ outliers	127900	1272 (2.56-2.52)

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	294	 89% 11%
1	B	294	 90% 8% ..
1	C	294	 91% 8% .
1	D	294	 88% 10% ..

## 2 Entry composition [i](#)

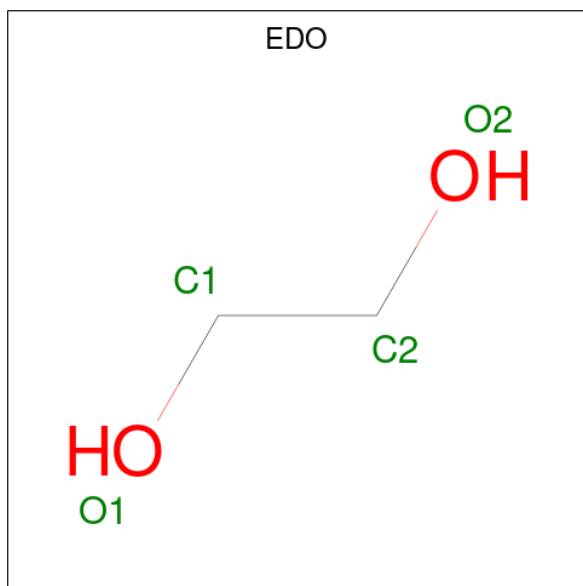
There are 5 unique types of molecules in this entry. The entry contains 9269 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Protein MEMO1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	293	Total 2300	C 1465	N 386	O 434	S 15	0	0	0
1	B	292	Total 2258	C 1437	N 376	O 430	S 15	0	0	0
1	C	293	Total 2277	C 1448	N 381	O 433	S 15	0	1	0
1	D	292	Total 2237	C 1428	N 364	O 430	S 15	0	0	0

- Molecule 2 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula: C<sub>2</sub>H<sub>6</sub>O<sub>2</sub>).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	C	O		
2	A	1	Total 4	C 2	O 2	0	0
2	A	1	Total 4	C 2	O 2	0	0

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Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	B	1	Total C O 4 2 2	0	0
2	B	1	Total C O 4 2 2	0	0
2	C	1	Total C O 4 2 2	0	0
2	C	1	Total C O 4 2 2	0	0
2	C	1	Total C O 4 2 2	0	0
2	C	1	Total C O 4 2 2	0	0
2	C	1	Total C O 4 2 2	0	0
2	C	1	Total C O 4 2 2	0	0
2	C	1	Total C O 4 2 2	0	0
2	C	1	Total C O 4 2 2	0	0
2	D	1	Total C O 4 2 2	0	0
2	D	1	Total C O 4 2 2	0	0
2	D	1	Total C O 4 2 2	0	0
2	D	1	Total C O 4 2 2	0	0

- Molecule 3 is COPPER (I) ION (three-letter code: CU1) (formula: Cu) (labeled as "Ligand of Interest" by depositor).

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

- Molecule 4 is GLYCEROL (three-letter code: GOL) (formula: C<sub>3</sub>H<sub>8</sub>O<sub>3</sub>).



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

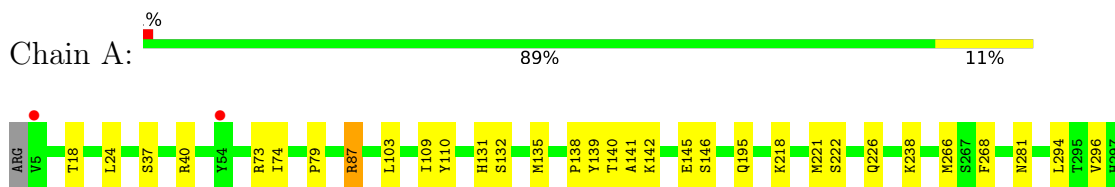
- Molecule 5 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	36	Total O 36 36	0	0
5	B	31	Total O 31 31	0	0
5	C	32	Total O 32 32	0	0
5	D	24	Total O 24 24	0	0

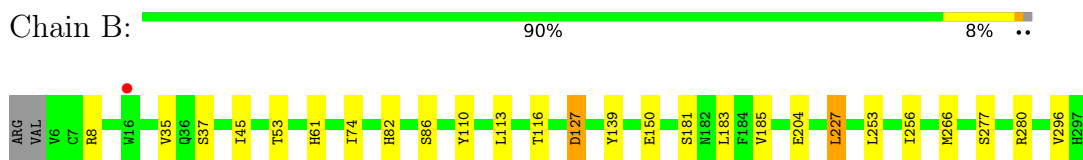
### 3 Residue-property plots [i](#)

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.

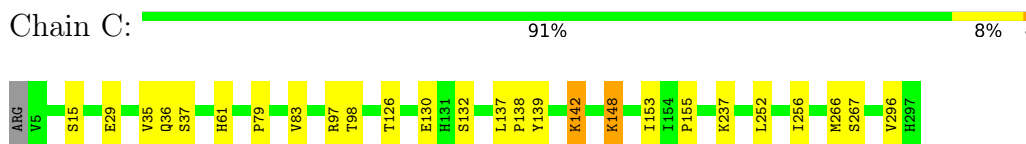
- Molecule 1: Protein MEMO1



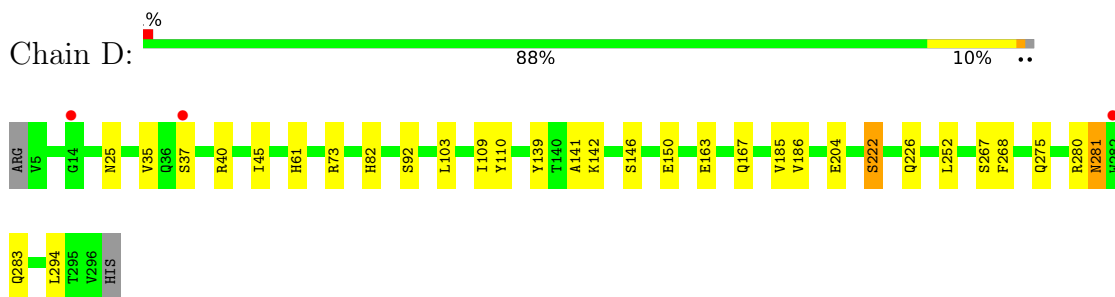
- Molecule 1: Protein MEMO1



- Molecule 1: Protein MEMO1



- Molecule 1: Protein MEMO1



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	140.63Å 87.16Å 98.59Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	42.91 – 2.55 47.85 – 2.55	Depositor EDS
% Data completeness (in resolution range)	99.9 (42.91-2.55) 91.9 (47.85-2.55)	Depositor EDS
$R_{merge}$	0.19	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.18 (at 2.54Å)	Xtrriage
Refinement program	PHENIX 1.19rc3_4028	Depositor
R, $R_{free}$	0.189 , 0.247 0.190 , 0.245	Depositor DCC
$R_{free}$ test set	2009 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	31.3	Xtrriage
Anisotropy	0.890	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.36 , 49.1	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.33$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	9269	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	44.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 12.41% 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 [i](#)

### 5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: CU1, GOL, EDO

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	0.26	0/2362	0.46	0/3205
1	B	0.27	0/2320	0.48	0/3156
1	C	0.27	0/2341	0.46	0/3181
1	D	0.25	0/2298	0.45	0/3127
All	All	0.26	0/9321	0.46	0/12669

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	2300	0	2167	16	0
1	B	2258	0	2077	9	0
1	C	2277	0	2119	13	0
1	D	2237	0	2055	17	0
2	A	8	0	12	0	0
2	B	8	0	12	0	0
2	C	32	0	48	2	0
2	D	16	0	24	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	B	6	0	8	0	0
5	A	36	0	0	1	0
5	B	31	0	0	0	0
5	C	32	0	0	0	0
5	D	24	0	0	1	0
All	All	9269	0	8522	52	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 (52) 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:D:37:SER:OG	1:D:40:ARG:HG3	1.80	0.81
1:B:82:HIS:HB3	1:D:82:HIS:HB3	1.65	0.78
1:D:25:ASN:OD1	1:D:142:LYS:NZ	2.21	0.73
1:D:281:ASN:HD22	1:D:283:GLN:H	1.35	0.72
1:A:73:ARG:HD3	1:A:109:ILE:HD11	1.81	0.63
1:B:8:ARG:NH2	1:B:127:ASP:O	2.32	0.62
1:D:73:ARG:HD3	1:D:109:ILE:HD11	1.81	0.61
1:B:35:VAL:HG11	1:B:61:HIS:HA	1.83	0.60
1:D:35:VAL:HG11	1:D:61:HIS:HA	1.84	0.59
1:D:275:GLN:NE2	5:D:401:HOH:O	2.35	0.59
1:D:281:ASN:ND2	1:D:283:GLN:H	2.02	0.57
1:C:252:LEU:O	1:C:256:ILE:HG13	2.06	0.56
1:C:130:GLU:HB3	2:C:302:EDO:H11	1.88	0.56
1:A:222:SER:O	1:A:226:GLN:HG2	2.07	0.54
1:B:113:LEU:O	1:B:116:THR:OG1	2.24	0.53
1:B:227:LEU:HD22	1:B:256:ILE:HG21	1.90	0.53
1:D:37:SER:HG	1:D:40:ARG:HG3	1.73	0.52
1:C:35:VAL:HG11	1:C:61:HIS:HA	1.92	0.52
1:A:266:MET:HG2	1:A:296:VAL:HG22	1.92	0.52
1:C:83:VAL:HG21	1:C:126:THR:HG23	1.93	0.51
1:B:74:ILE:HG12	1:B:183:LEU:HD23	1.93	0.50
1:D:186:VAL:HG21	1:D:252:LEU:HD22	1.93	0.50
1:B:266:MET:HG2	1:B:296:VAL:HG22	1.94	0.50
1:A:24:LEU:HD21	1:A:138:PRO:HB3	1.94	0.50
1:C:266:MET:HG2	1:C:296:VAL:HG22	1.94	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:45:ILE:HD11	1:B:185:VAL:HG22	1.93	0.49
1:A:268:PHE:HD2	1:A:294:LEU:HB2	1.77	0.49
1:A:37:SER:HB2	1:A:40:ARG:HG3	1.95	0.48
1:C:29:GLU:OE2	1:C:142:LYS:NZ	2.35	0.48
1:D:73:ARG:HG3	1:D:73:ARG:HH11	1.80	0.47
1:D:222:SER:O	1:D:226:GLN:HG2	2.15	0.46
1:A:195:GLN:HG2	1:C:97:ARG:HH12	1.81	0.46
1:A:103:LEU:HD21	1:A:141:ALA:HB2	1.97	0.45
1:D:45:ILE:HD11	1:D:185:VAL:HG22	1.97	0.45
1:B:227:LEU:HA	1:B:253:LEU:HD22	1.98	0.45
1:D:163:GLU:O	1:D:167:GLN:HG3	2.17	0.44
1:D:268:PHE:HD2	1:D:294:LEU:HB2	1.82	0.43
1:C:36:GLN:H	2:C:305:EDO:H21	1.83	0.43
1:A:18:THR:HG23	1:A:24:LEU:HB2	2.01	0.43
1:C:237:LYS:HG2	1:D:204:GLU:HB3	2.01	0.42
1:D:103:LEU:HD21	1:D:141:ALA:HB2	2.02	0.42
1:C:153:ILE:HD12	1:C:155:PRO:HG3	2.02	0.42
1:C:148:LYS:HB2	1:C:148:LYS:HE2	1.38	0.41
1:A:74:ILE:HD12	1:A:140:THR:HG23	2.02	0.41
1:C:79:PRO:HD2	1:C:132:SER:HB2	2.03	0.41
1:C:137:LEU:HB2	1:C:138:PRO:HD3	2.02	0.41
1:A:131:HIS:O	1:A:135:MET:HG2	2.21	0.41
1:A:218:LYS:HA	1:A:221:MET:HB2	2.03	0.41
1:A:79:PRO:HD2	1:A:132:SER:HB2	2.03	0.40
1:A:142:LYS:HD2	1:A:142:LYS:HA	1.91	0.40
1:A:87:ARG:HD2	5:A:418:HOH:O	2.21	0.40
1:A:142:LYS:HE3	1:A:145:GLU:OE2	2.21	0.40

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	291/294 (99%)	284 (98%)	7 (2%)	0	100	100
1	B	290/294 (99%)	282 (97%)	8 (3%)	0	100	100
1	C	292/294 (99%)	286 (98%)	6 (2%)	0	100	100
1	D	290/294 (99%)	282 (97%)	8 (3%)	0	100	100
All	All	1163/1176 (99%)	1134 (98%)	29 (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	242/256 (94%)	236 (98%)	6 (2%)	47	62
1	B	231/256 (90%)	219 (95%)	12 (5%)	23	30
1	C	236/256 (92%)	229 (97%)	7 (3%)	41	55
1	D	228/256 (89%)	219 (96%)	9 (4%)	32	44
All	All	937/1024 (92%)	903 (96%)	34 (4%)	35	47

All (34) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	87	ARG
1	A	110	TYR
1	A	139	TYR
1	A	146	SER
1	A	238	LYS
1	A	281	ASN
1	B	37	SER
1	B	53	THR
1	B	86	SER
1	B	110	TYR
1	B	127	ASP
1	B	139	TYR
1	B	150	GLU

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Mol	Chain	Res	Type
1	B	181	SER
1	B	204	GLU
1	B	227	LEU
1	B	277	SER
1	B	280	ARG
1	C	15	SER
1	C	37	SER
1	C	98	THR
1	C	139	TYR
1	C	142	LYS
1	C	148	LYS
1	C	267	SER
1	D	92	SER
1	D	110	TYR
1	D	139	TYR
1	D	146	SER
1	D	150	GLU
1	D	222	SER
1	D	267	SER
1	D	280	ARG
1	D	281	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	27	GLN
1	B	131	HIS
1	C	262	ASN
1	D	262	ASN
1	D	281	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 21 ligands modelled in this entry, 4 are monoatomic - leaving 17 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	$\# Z  > 2$
2	EDO	D	304	-	3,3,3	0.46	0	2,2,2	0.26	0
2	EDO	B	303	-	3,3,3	0.46	0	2,2,2	0.38	0
2	EDO	C	305	-	3,3,3	0.45	0	2,2,2	0.37	0
2	EDO	C	308	-	3,3,3	0.46	0	2,2,2	0.31	0
2	EDO	B	302	-	3,3,3	0.44	0	2,2,2	0.38	0
2	EDO	C	307	-	3,3,3	0.46	0	2,2,2	0.34	0
4	GOL	B	301	-	5,5,5	0.84	0	5,5,5	1.02	0
2	EDO	D	303	-	3,3,3	0.45	0	2,2,2	0.35	0
2	EDO	D	302	-	3,3,3	0.48	0	2,2,2	0.23	0
2	EDO	C	304	-	3,3,3	0.47	0	2,2,2	0.36	0
2	EDO	A	302	-	3,3,3	0.45	0	2,2,2	0.36	0
2	EDO	C	306	-	3,3,3	0.47	0	2,2,2	0.34	0
2	EDO	D	301	-	3,3,3	0.48	0	2,2,2	0.30	0
2	EDO	C	302	-	3,3,3	0.43	0	2,2,2	0.32	0
2	EDO	A	301	-	3,3,3	0.47	0	2,2,2	0.39	0
2	EDO	C	303	-	3,3,3	0.45	0	2,2,2	0.36	0
2	EDO	C	301	-	3,3,3	0.46	0	2,2,2	0.37	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
2	EDO	D	304	-	-	0/1/1/1	-
2	EDO	B	303	-	-	0/1/1/1	-
2	EDO	C	305	-	-	0/1/1/1	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	EDO	C	308	-	-	0/1/1/1	-
2	EDO	B	302	-	-	0/1/1/1	-
2	EDO	C	307	-	-	1/1/1/1	-
4	GOL	B	301	-	-	2/4/4/4	-
2	EDO	D	303	-	-	0/1/1/1	-
2	EDO	D	302	-	-	0/1/1/1	-
2	EDO	C	304	-	-	0/1/1/1	-
2	EDO	A	302	-	-	0/1/1/1	-
2	EDO	C	306	-	-	0/1/1/1	-
2	EDO	D	301	-	-	0/1/1/1	-
2	EDO	C	302	-	-	0/1/1/1	-
2	EDO	A	301	-	-	0/1/1/1	-
2	EDO	C	303	-	-	1/1/1/1	-
2	EDO	C	301	-	-	0/1/1/1	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (4) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	B	301	GOL	O1-C1-C2-C3
4	B	301	GOL	O1-C1-C2-O2
2	C	303	EDO	O1-C1-C2-O2
2	C	307	EDO	O1-C1-C2-O2

There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	305	EDO	1	0
2	C	302	EDO	1	0

## 5.7 Other polymers [i](#)

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	293/294 (99%)	-0.06	2 (0%) 87 90	25, 39, 57, 88	0
1	B	292/294 (99%)	0.06	1 (0%) 94 96	29, 46, 75, 89	0
1	C	293/294 (99%)	-0.02	0 100 100	27, 41, 57, 76	0
1	D	292/294 (99%)	0.10	3 (1%) 82 86	28, 47, 76, 85	0
All	All	1170/1176 (99%)	0.02	6 (0%) 91 94	25, 43, 68, 89	0

All (6) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	5	VAL	5.0
1	A	54	TYR	2.2
1	D	14	GLY	2.2
1	D	282	TRP	2.1
1	D	37	SER	2.1
1	B	16	TRP	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,



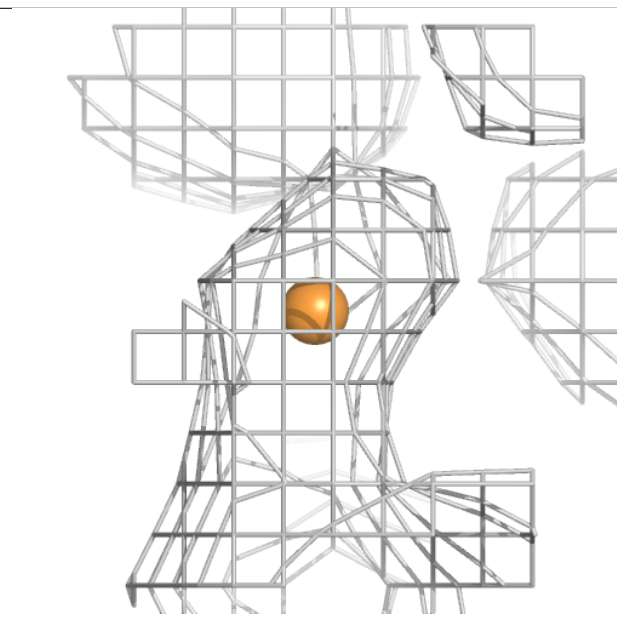
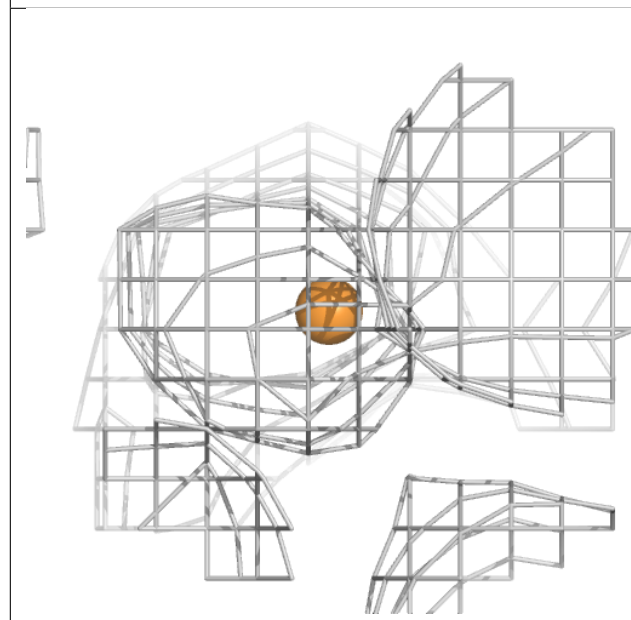
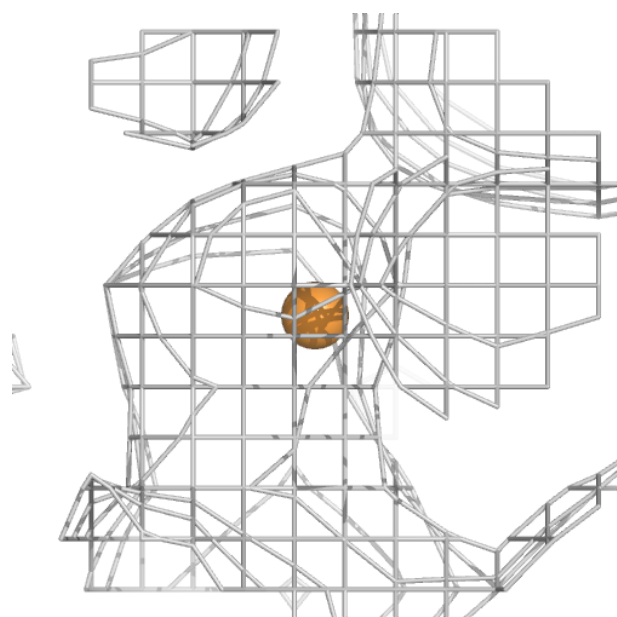
median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å <sup>2</sup> )	Q<0.9
3	CU1	D	305	1/1	0.63	0.30	119,119,119,119	0
3	CU1	C	309	1/1	0.65	0.17	73,73,73,73	1
2	EDO	D	302	4/4	0.67	0.36	45,47,50,55	0
2	EDO	C	308	4/4	0.69	0.30	44,51,57,58	0
2	EDO	B	303	4/4	0.69	0.29	52,66,66,69	0
2	EDO	C	305	4/4	0.73	0.27	61,62,64,66	0
2	EDO	D	304	4/4	0.76	0.22	43,45,47,53	0
2	EDO	C	306	4/4	0.78	0.20	45,46,54,55	0
2	EDO	D	301	4/4	0.78	0.26	45,48,50,54	0
3	CU1	B	304	1/1	0.81	0.31	60,60,60,60	1
3	CU1	A	303	1/1	0.85	0.36	67,67,67,67	1
4	GOL	B	301	6/6	0.85	0.23	39,45,53,56	0
2	EDO	A	301	4/4	0.86	0.16	46,56,60,62	0
2	EDO	C	301	4/4	0.88	0.25	43,45,56,56	0
2	EDO	C	307	4/4	0.90	0.16	35,36,40,41	0
2	EDO	D	303	4/4	0.90	0.15	58,63,65,71	0
2	EDO	B	302	4/4	0.90	0.20	50,51,58,63	0
2	EDO	C	303	4/4	0.90	0.23	41,44,48,57	0
2	EDO	C	304	4/4	0.91	0.18	44,45,46,48	0
2	EDO	C	302	4/4	0.93	0.49	38,39,41,42	0
2	EDO	A	302	4/4	0.94	0.15	48,55,59,61	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.

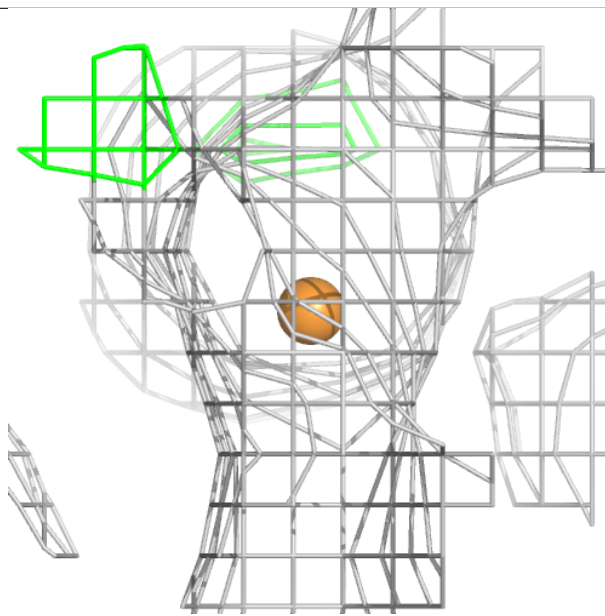
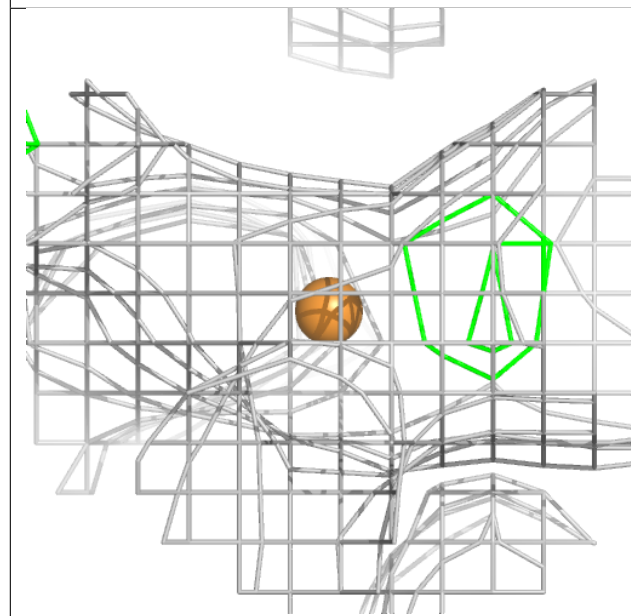
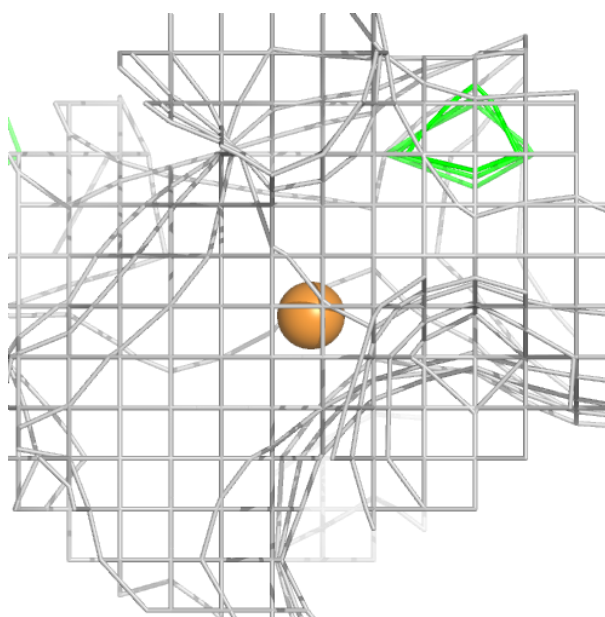
**Electron density around CU1 D 305:**

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



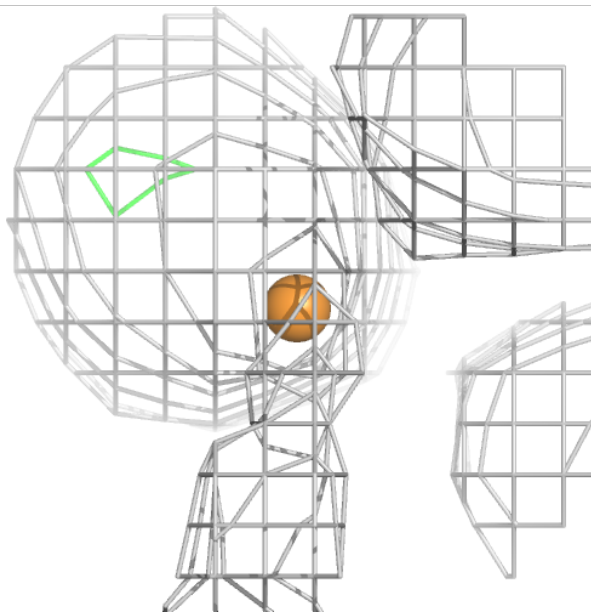
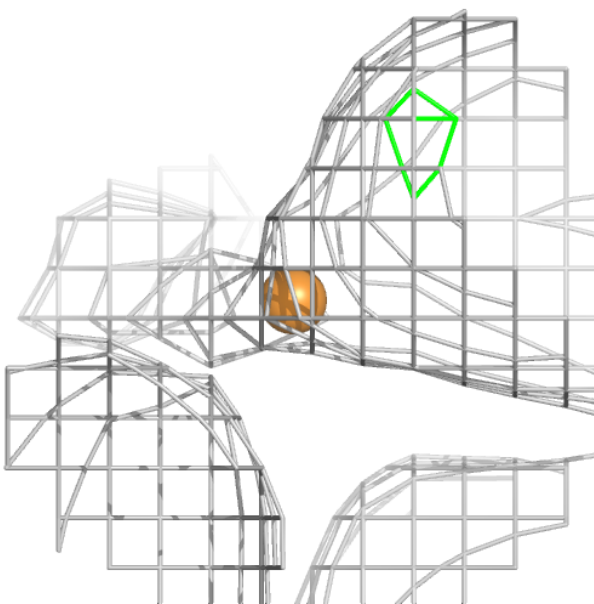
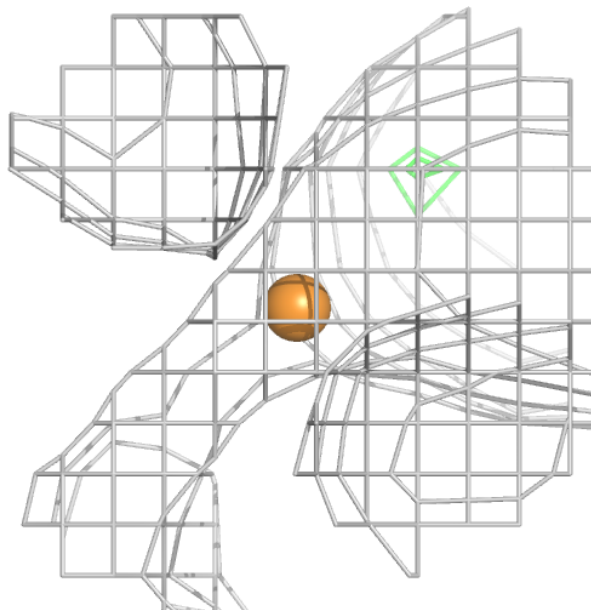
**Electron density around CU1 C 309:**

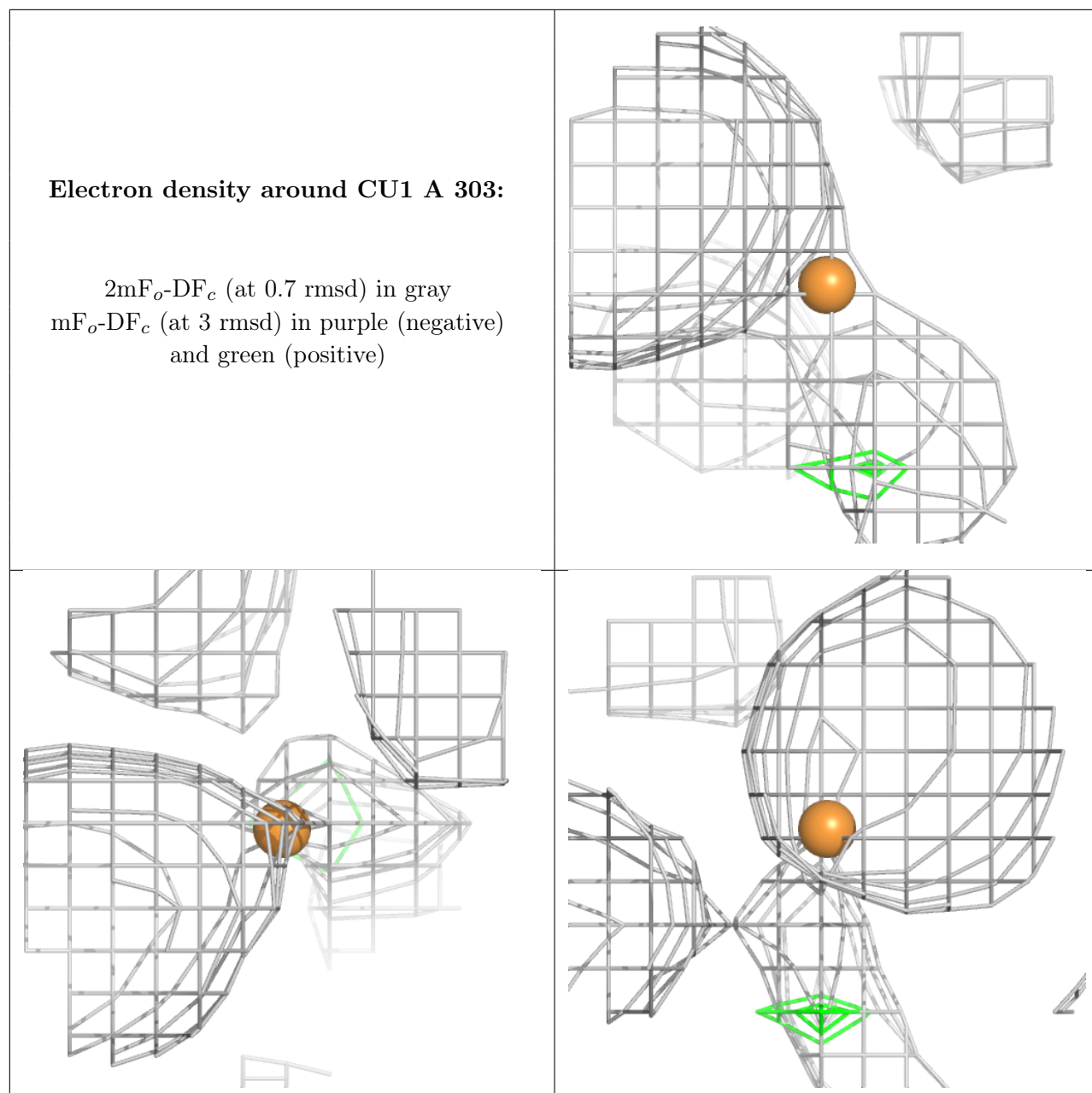
$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)



**Electron density around CU1 B 304:**

$2mF_o-DF_c$  (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.