



# Full wwPDB X-ray Structure Validation Report ⓘ

Nov 19, 2023 – 05:29 PM JST

PDB ID : 6LTU  
Title : Crystal structure of Cas12i2 ternary complex with double Mg<sup>2+</sup> bound in catalytic pocket  
Authors : Huang, X.; Sun, W.; Cheng, Z.; Chen, M.; Li, X.; Wang, J.; Sheng, G.; Gong, W.; Wang, Y.  
Deposited on : 2020-01-23  
Resolution : 2.57 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)  
Xtrriage (Phenix) : 1.13  
EDS : 2.36  
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) : 2.36

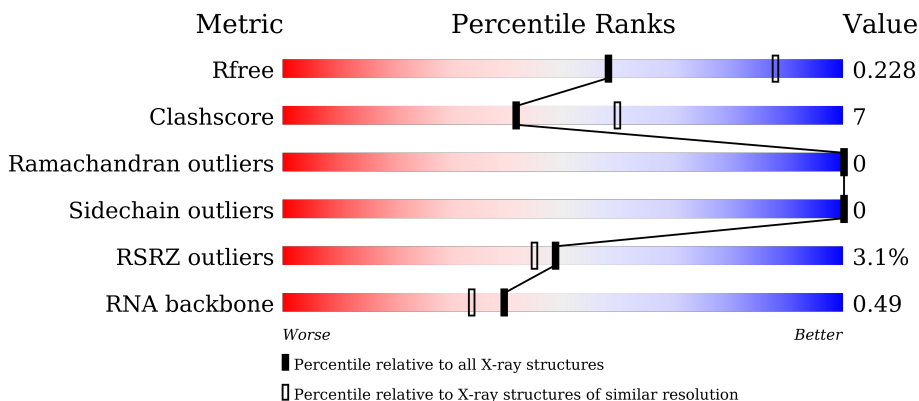
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.57 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	3676 (2.60-2.56)
Clashscore	141614	4049 (2.60-2.56)
Ramachandran outliers	138981	3979 (2.60-2.56)
Sidechain outliers	138945	3979 (2.60-2.56)
RSRZ outliers	127900	3614 (2.60-2.56)
RNA backbone	3102	1075 (2.90-2.26)

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	1055	 3% 83% 15%
2	B	58	 3% 47% 21% 16% 16%
3	D	12	 8% 42% 42% 8% 8%

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Mol	Chain	Length	Quality of chain
4	C	35	
5	E	5	

## 2 Entry composition

There are 8 unique types of molecules in this entry. The entry contains 10704 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 Cas12i2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	1034	8312	5280	1467	1538	27	0	1	0

- Molecule 2 is a RNA chain called RNA (56-mer).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	P			
2	B	49	1045	469	191	337	48	0	0	0

- Molecule 3 is a DNA chain called DNA (5'-D(\*GP\*CP\*CP\*GP\*CP\*TP\*TP\*TP\*CP\*TP\*T)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	P			
3	D	11	217	106	32	69	10	0	0	0

- Molecule 4 is a DNA chain called DNA (35-MER).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	P			
4	C	35	715	341	127	212	35	0	0	0

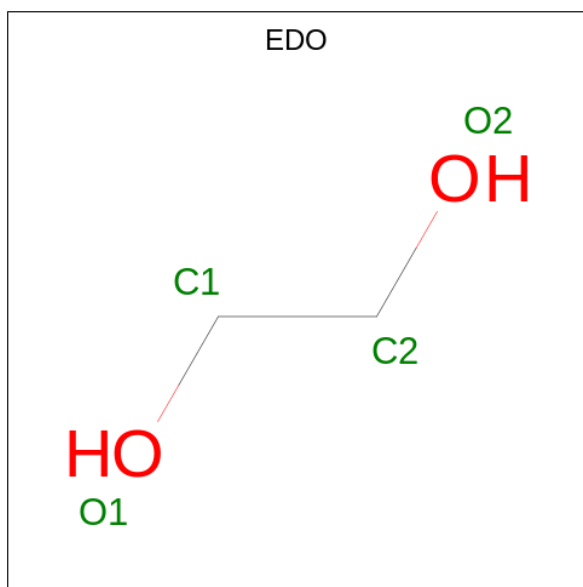
- Molecule 5 is a DNA chain called trans ssDNA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	P			
5	E	5	83	38	13	27	5	0	0	0

- Molecule 6 is MAGNESIUM ION (three-letter code: MG) (formula: Mg) (labeled as "Ligand of Interest" by depositor).

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

- Molecule 7 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
7	A	1	Total C O 4 2 2	0	0

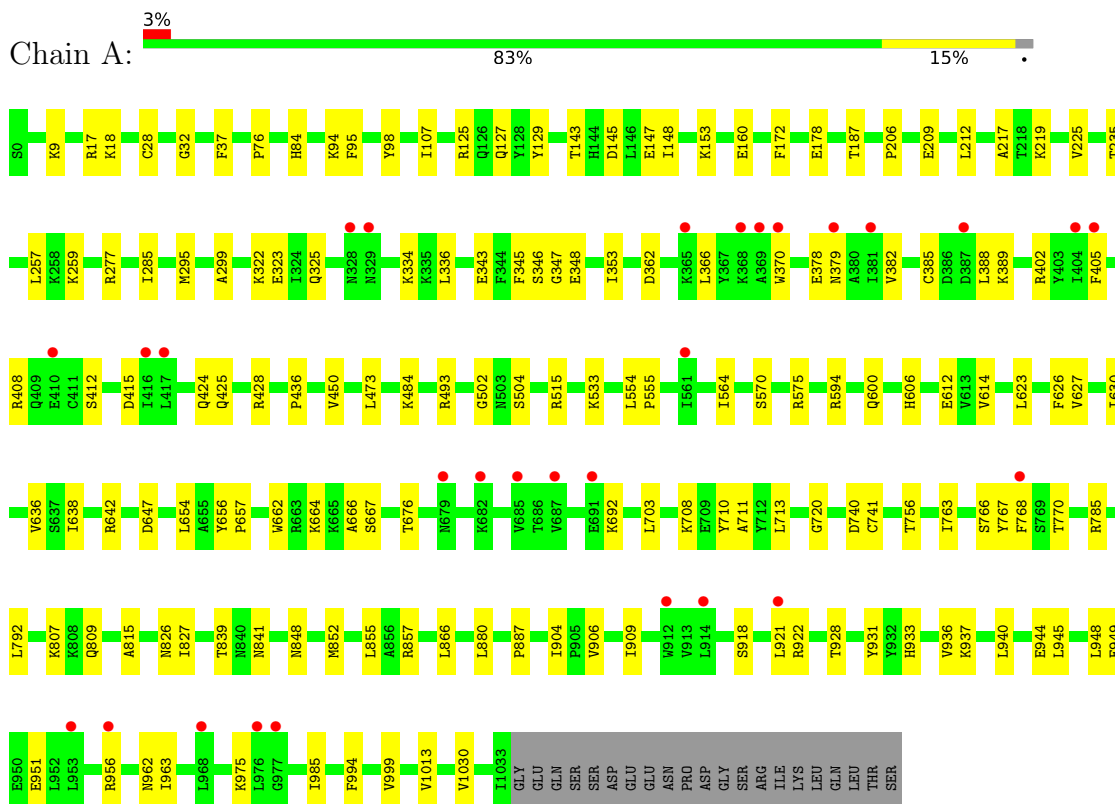
- Molecule 8 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
8	A	229	Total O 229 229	0	0
8	B	48	Total O 48 48	0	0
8	D	16	Total O 16 16	0	0
8	C	31	Total O 31 31	0	0
8	E	1	Total O 1 1	0	0

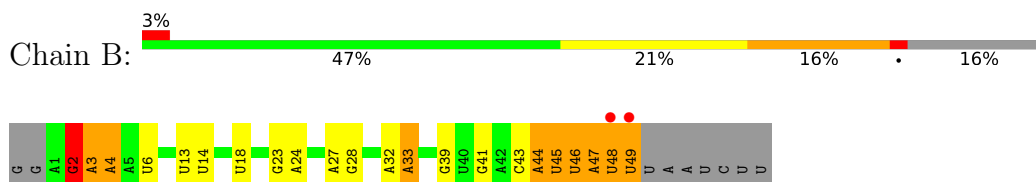
### 3 Residue-property plots [i](#)

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.

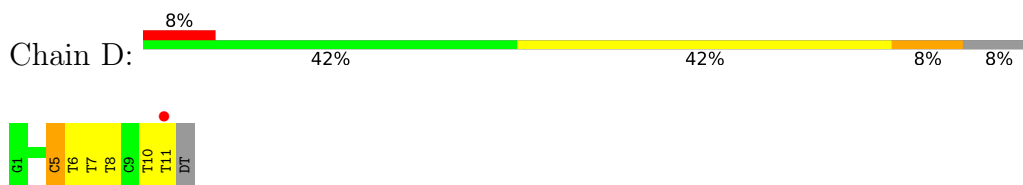
- Molecule 1: Cas12i2



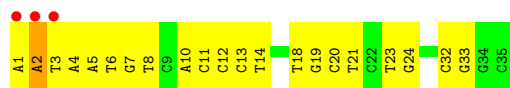
- Molecule 2: RNA (56-mer)



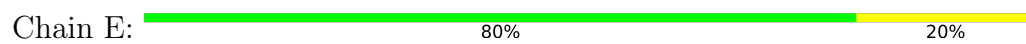
- Molecule 3: DNA (5'-D(\*GP\*CP\*CP\*GP\*CP\*TP\*TP\*TP\*CP\*TP\*T)-3')



## ● Molecule 4: DNA (35-MER)



## ● Molecule 5: trans ssDNA



## 4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	94.33Å 123.58Å 281.22Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	46.87 – 2.57 46.87 – 2.57	Depositor EDS
% Data completeness (in resolution range)	97.0 (46.87-2.57) 97.0 (46.87-2.57)	Depositor EDS
$R_{merge}$	0.16	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.76 (at 2.58Å)	Xtrriage
Refinement program	PHENIX 1.18.2_3874	Depositor
R, $R_{free}$	0.202 , 0.229 0.202 , 0.228	Depositor DCC
$R_{free}$ test set	2504 reflections (4.90%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	41.7	Xtrriage
Anisotropy	0.205	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.30 , 37.3	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.48$ , $\langle L^2 \rangle = 0.32$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
$F_o, F_c$ correlation	0.93	EDS
Total number of atoms	10704	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	49.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.98% 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: MG, 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	0.42	0/8481	0.55	0/11451
2	B	0.38	0/1170	1.06	10/1822 (0.5%)
3	D	0.66	0/240	1.09	1/368 (0.3%)
4	C	0.61	0/800	1.15	4/1232 (0.3%)
5	E	0.68	0/91	1.09	0/138
All	All	0.44	0/10782	0.72	15/15011 (0.1%)

There are no bond length outliers.

All (15) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	3	A	C4'-C3'-O3'	-10.97	86.36	109.40
2	B	47	A	N9-C1'-C2'	-8.68	102.45	112.00
4	C	3	DT	O5'-P-OP2	-8.68	97.89	105.70
2	B	47	A	C1'-C2'-O2'	-8.23	85.90	110.60
4	C	2	DA	C4-N9-C1'	-7.91	112.06	126.30
4	C	2	DA	C8-N9-C1'	7.89	141.91	127.70
2	B	3	A	O5'-P-OP1	-7.41	99.03	105.70
2	B	2	G	C4'-C3'-O3'	-7.41	93.85	109.40
2	B	47	A	C3'-C2'-C1'	7.19	107.25	101.50
2	B	47	A	C4'-C3'-O3'	6.64	126.29	113.00
2	B	47	A	O4'-C1'-N9	6.12	113.09	108.20
2	B	48	U	C1'-O4'-C4'	5.68	114.45	109.90
3	D	5	DC	O4'-C4'-C3'	-5.46	102.32	104.50
4	C	2	DA	P-O3'-C3'	5.41	126.19	119.70
2	B	4	A	C1'-C2'-O2'	-5.13	95.21	110.60

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	8312	0	8277	112	0
2	B	1045	0	529	15	0
3	D	217	0	128	7	0
4	C	715	0	396	16	0
5	E	83	0	45	1	0
6	A	2	0	0	0	0
6	B	1	0	0	0	0
7	A	4	0	6	1	0
8	A	229	0	0	3	0
8	B	48	0	0	0	0
8	C	31	0	0	0	0
8	D	16	0	0	0	0
8	E	1	0	0	0	0
All	All	10704	0	9381	138	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 7.

All (138) 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:428:ARG:NH2	2:B:32:A:OP1	2.08	0.87
4:C:1:DA:H2'	4:C:1:DA:N3	1.90	0.86
1:A:259:LYS:HD3	3:D:10:DT:H2''	1.59	0.83
1:A:921:LEU:O	1:A:921:LEU:HD23	1.85	0.75
4:C:2:DA:H3'	4:C:2:DA:OP2	1.87	0.74
1:A:178:GLU:HB3	3:D:8:DT:H5''	1.71	0.72
1:A:807:LYS:HG2	1:A:855:LEU:HD11	1.69	0.72
1:A:918:SER:CB	1:A:922:ARG:HH12	2.03	0.71
1:A:994:PHE:HB3	1:A:1013:VAL:HG21	1.72	0.69
4:C:2:DA:H3'	4:C:2:DA:P	2.34	0.67
1:A:956:ARG:O	1:A:962:ASN:ND2	2.20	0.67
1:A:336:LEU:HD21	1:A:353:ILE:HG21	1.78	0.66
1:A:370:TRP:CD2	1:A:405:PHE:HD1	2.15	0.64
1:A:172:PHE:O	1:A:277:ARG:NH1	2.31	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:343:GLU:HA	1:A:343:GLU:OE2	1.97	0.62
1:A:259:LYS:HD3	3:D:10:DT:C2'	2.30	0.61
1:A:951:GLU:HB3	1:A:963:ILE:HD11	1.81	0.61
1:A:945:LEU:HB3	1:A:948:LEU:HD12	1.82	0.61
1:A:933:HIS:O	1:A:937:LYS:HG2	2.01	0.61
1:A:370:TRP:CZ2	1:A:405:PHE:HB2	2.37	0.60
1:A:570:SER:HB3	7:A:1103:EDO:H12	1.85	0.57
1:A:370:TRP:CE3	1:A:408:ARG:HD2	2.41	0.56
1:A:614:VAL:O	1:A:626:PHE:HB2	2.05	0.56
1:A:18:LYS:NZ	2:B:2:G:N7	2.51	0.55
1:A:654:LEU:HD12	1:A:703:LEU:HD22	1.88	0.55
1:A:235:THR:HG23	3:D:10:DT:OP1	2.05	0.55
1:A:370:TRP:CE2	1:A:405:PHE:HD1	2.23	0.55
1:A:366:LEU:HD21	1:A:370:TRP:CZ2	2.41	0.55
1:A:692:LYS:HG2	1:A:713:LEU:HD13	1.88	0.54
2:B:45:U:H2'	2:B:46:U:C6	2.42	0.54
1:A:379:ASN:HA	1:A:382:VAL:HG22	1.90	0.54
1:A:928:THR:HA	1:A:931:TYR:CE2	2.43	0.54
1:A:904:ILE:HD12	1:A:985:ILE:HG13	1.89	0.53
2:B:44:A:H2'	2:B:45:U:C6	2.43	0.53
2:B:48:U:H2'	2:B:49:U:C6	2.42	0.53
1:A:708:LYS:HG3	1:A:756:THR:HA	1.91	0.53
1:A:662:TRP:CZ3	1:A:740:ASP:HB2	2.44	0.52
1:A:921:LEU:HD23	1:A:921:LEU:C	2.29	0.52
1:A:921:LEU:HD11	1:A:949:GLU:HG3	1.90	0.52
1:A:928:THR:HA	1:A:931:TYR:CD2	2.44	0.52
1:A:600:GLN:HA	1:A:606:HIS:CD2	2.45	0.52
3:D:7:DT:H2'	3:D:8:DT:C6	2.45	0.52
1:A:918:SER:C	1:A:922:ARG:NH1	2.64	0.51
1:A:187:THR:HG23	1:A:257:LEU:HD22	1.92	0.51
1:A:612:GLU:HB2	1:A:630:ILE:HD11	1.92	0.51
1:A:944:GLU:HB2	1:A:975:LYS:NZ	2.25	0.51
1:A:711:ALA:HA	1:A:763:ILE:HD11	1.94	0.50
1:A:918:SER:O	1:A:922:ARG:HG2	2.11	0.50
1:A:17:ARG:NH1	8:A:1203:HOH:O	2.33	0.50
1:A:623:LEU:HD12	1:A:627:VAL:HG11	1.94	0.50
1:A:594:ARG:HB3	1:A:827:ILE:HD13	1.94	0.49
1:A:918:SER:CB	1:A:922:ARG:NH1	2.74	0.49
1:A:378:GLU:OE2	1:A:402:ARG:NE	2.43	0.49
1:A:493:ARG:NH1	8:A:1207:HOH:O	2.42	0.49
1:A:212:LEU:HD11	1:A:219:LYS:HG3	1.94	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:841:ASN:N	2:B:33:A:OP1	2.46	0.49
1:A:322:LYS:HA	1:A:325:GLN:HG2	1.93	0.49
4:C:18:DT:H2'	4:C:19:DG:C8	2.48	0.49
1:A:385:CYS:O	1:A:389:LYS:HG3	2.13	0.49
2:B:49:U:OP2	2:B:49:U:H6	1.96	0.49
1:A:346:SER:OG	2:B:41:G:H5''	2.13	0.48
1:A:504:SER:HB3	1:A:555:PRO:HA	1.95	0.48
1:A:370:TRP:CH2	1:A:405:PHE:HB2	2.48	0.48
1:A:656:TYR:HB3	1:A:657:PRO:HD3	1.95	0.48
1:A:143:THR:O	1:A:147:GLU:HG3	2.13	0.48
1:A:450:VAL:HB	1:A:564:ILE:HB	1.95	0.48
1:A:944:GLU:HB2	1:A:975:LYS:HZ3	1.79	0.47
1:A:815:ALA:HB1	1:A:866:LEU:HD12	1.95	0.47
1:A:906:VAL:HA	1:A:909:ILE:HD12	1.97	0.47
1:A:378:GLU:O	1:A:382:VAL:HG13	2.14	0.47
1:A:975:LYS:HE3	1:A:975:LYS:HB2	1.53	0.46
1:A:37:PHE:HA	1:A:299:ALA:HB1	1.98	0.46
1:A:666:ALA:HB1	1:A:710:TYR:CE1	2.51	0.46
1:A:153:LYS:HE3	1:A:160:GLU:CG	2.46	0.45
1:A:379:ASN:O	1:A:382:VAL:HG22	2.16	0.45
1:A:515:ARG:HD3	2:B:23:G:OP1	2.16	0.45
4:C:8:DT:H6	4:C:8:DT:H5''	1.81	0.45
4:C:1:DA:N3	4:C:1:DA:C2'	2.71	0.45
1:A:600:GLN:HG2	1:A:606:HIS:CD2	2.52	0.45
1:A:84:HIS:CE1	1:A:127:GLN:HA	2.52	0.45
1:A:94:LYS:HG3	1:A:98:TYR:CE2	2.51	0.45
1:A:502:GLY:HA2	1:A:554:LEU:HD23	1.99	0.44
4:C:12:DC:H2'	4:C:13:DC:C6	2.52	0.44
1:A:623:LEU:HD23	1:A:1030:VAL:CG1	2.47	0.44
1:A:28:CYS:SG	1:A:285:ILE:HG21	2.57	0.44
1:A:642:ARG:HD2	2:B:14:U:O2	2.17	0.44
1:A:412:SER:OG	1:A:415:ASP:OD1	2.35	0.44
1:A:424:GLN:O	1:A:428:ARG:HG3	2.17	0.44
2:B:27:A:H2'	2:B:28:G:C8	2.52	0.44
1:A:767:TYR:CE2	1:A:792:LEU:HD23	2.53	0.43
4:C:6:DT:H2''	4:C:7:DG:O4'	2.17	0.43
1:A:388:LEU:HA	1:A:388:LEU:HD23	1.76	0.43
1:A:839:THR:HB	2:B:33:A:H4'	2.00	0.43
3:D:11:DT:H6	3:D:11:DT:H2'	1.41	0.43
2:B:43:C:H2'	2:B:44:A:C8	2.54	0.43
1:A:362:ASP:OD1	1:A:388:LEU:HD21	2.19	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:217:ALA:HB2	1:A:225:VAL:HG21	2.00	0.43
1:A:809:GLN:HE21	1:A:809:GLN:HB2	1.61	0.43
4:C:20:DC:C6	4:C:21:DT:H72	2.54	0.43
4:C:23:DT:H2''	4:C:24:DG:H5''	2.00	0.42
1:A:125:ARG:HD2	1:A:129:TYR:CZ	2.54	0.42
1:A:206:PRO:O	1:A:209:GLU:HB3	2.19	0.42
1:A:921:LEU:C	1:A:921:LEU:CD2	2.88	0.42
4:C:2:DA:OP2	4:C:2:DA:H2'	2.19	0.42
1:A:766:SER:O	1:A:770:THR:HG23	2.19	0.42
2:B:45:U:H2'	2:B:46:U:H6	1.83	0.42
2:B:45:U:H2'	2:B:46:U:O4'	2.19	0.42
1:A:32:GLY:HA3	1:A:295:MET:SD	2.60	0.42
1:A:848:ASN:O	1:A:852:MET:HG3	2.18	0.42
1:A:323:GLU:OE1	1:A:425:GLN:NE2	2.51	0.42
1:A:9:LYS:HG3	1:A:575:ARG:HG2	2.01	0.42
4:C:32:DC:H2'	4:C:33:DG:C8	2.54	0.42
1:A:277:ARG:NH2	8:A:1214:HOH:O	2.45	0.41
1:A:662:TRP:CE3	1:A:740:ASP:HB2	2.55	0.41
1:A:887:PRO:HG2	1:A:999:VAL:HG21	2.01	0.41
1:A:76:PRO:O	1:A:436:PRO:HA	2.21	0.41
1:A:880:LEU:HA	5:E:22:DA:H1'	2.03	0.41
4:C:13:DC:H2'	4:C:14:DT:C6	2.55	0.41
1:A:533:LYS:HA	1:A:533:LYS:HD2	1.86	0.41
3:D:5:DC:H2'	3:D:6:DT:C6	2.55	0.41
4:C:4:DA:H2''	4:C:5:DA:H8	1.86	0.41
1:A:145:ASP:HA	1:A:148:ILE:HD12	2.03	0.41
1:A:664:LYS:O	1:A:667:SER:HB3	2.21	0.41
1:A:703:LEU:HD21	1:A:741:CYS:HB3	2.03	0.41
1:A:768:PHE:CG	1:A:785:ARG:HG2	2.56	0.41
1:A:676:THR:HB	1:A:720:GLY:HA3	2.03	0.41
1:A:826:ASN:O	1:A:826:ASN:ND2	2.53	0.41
1:A:345:PHE:CE1	1:A:347:GLY:HA3	2.55	0.41
1:A:473:LEU:O	1:A:484:LYS:HA	2.21	0.41
4:C:10:DA:H2'	4:C:11:DC:C6	2.56	0.41
1:A:95:PHE:CZ	1:A:107:ILE:HD13	2.56	0.40
1:A:636:VAL:HG12	1:A:638:ILE:HG12	2.03	0.40
1:A:636:VAL:HG13	1:A:647:ASP:OD1	2.21	0.40
1:A:857:ARG:HB2	4:C:20:DC:H5''	2.03	0.40
1:A:921:LEU:HB2	1:A:936:VAL:HG11	2.03	0.40
1:A:940:LEU:HD13	1:A:949:GLU:HB2	2.02	0.40
1:A:334:LYS:HA	1:A:334:LYS:HD2	1.66	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:348:GLU:H	1:A:348:GLU:HG3	1.64	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	1033/1055 (98%)	1010 (98%)	23 (2%)	0	<b>100</b>   <b>100</b>

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	890/931 (96%)	890 (100%)	0	<b>100</b>   <b>100</b>

There are no protein residues with a non-rotameric sidechain to report.

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

Mol	Chain	Res	Type
1	A	164	ASN
1	A	641	ASN

### 5.3.3 RNA [i](#)

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
2	B	48/58 (82%)	13 (27%)	3 (6%)

All (13) RNA backbone outliers are listed below:

Mol	Chain	Res	Type
2	B	2	G
2	B	3	A
2	B	4	A
2	B	6	U
2	B	13	U
2	B	18	U
2	B	24	A
2	B	33	A
2	B	39	G
2	B	44	A
2	B	46	U
2	B	47	A
2	B	49	U

All (3) RNA pucker outliers are listed below:

Mol	Chain	Res	Type
2	B	2	G
2	B	3	A
2	B	45	U

## 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 4 ligands modelled in this entry, 3 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	$\# Z  > 2$
7	EDO	A	1103	-	3,3,3	0.46	0	2,2,2	0.36	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
7	EDO	A	1103	-	-	0/1/1/1	-

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.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
7	A	1103	EDO	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

### 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, 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	1034/1055 (98%)	0.05	29 (2%) 53 49	26, 45, 85, 107	0
2	B	49/58 (84%)	-0.50	2 (4%) 37 33	32, 40, 108, 141	0
3	D	11/12 (91%)	0.01	1 (9%) 9 7	36, 39, 62, 97	0
4	C	35/35 (100%)	-0.01	3 (8%) 10 8	35, 45, 129, 140	0
5	E	5/5 (100%)	-0.03	0 100 100	46, 49, 59, 69	0
All	All	1134/1165 (97%)	0.03	35 (3%) 49 45	26, 45, 87, 141	0

All (35) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	679	ASN	6.2
1	A	328	ASN	4.3
1	A	329	ASN	4.2
1	A	921	LEU	4.1
2	B	49	U	4.0
1	A	682	LYS	3.9
4	C	1	DA	3.7
1	A	977	GLY	3.6
4	C	2	DA	3.5
1	A	953	LEU	3.3
1	A	687	VAL	3.0
2	B	48	U	2.9
1	A	370	TRP	2.8
1	A	404	ILE	2.7
1	A	368	LYS	2.7
3	D	11	DT	2.6
1	A	956	ARG	2.6
1	A	387	ASP	2.6
1	A	912	TRP	2.4
1	A	369	ALA	2.4

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Mol	Chain	Res	Type	RSRZ
1	A	365	LYS	2.4
1	A	417	LEU	2.4
1	A	914	LEU	2.3
1	A	691	GLU	2.3
1	A	416	ILE	2.3
1	A	968	LEU	2.2
1	A	410	GLU	2.2
1	A	768	PHE	2.2
1	A	561	ILE	2.2
1	A	685	VAL	2.1
1	A	976	LEU	2.1
1	A	379	ASN	2.1
1	A	381	ILE	2.1
1	A	405	PHE	2.1
4	C	3	DT	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 [i](#)

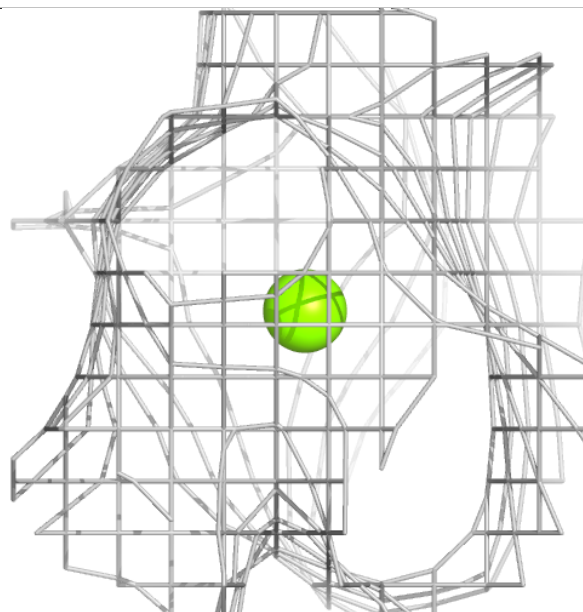
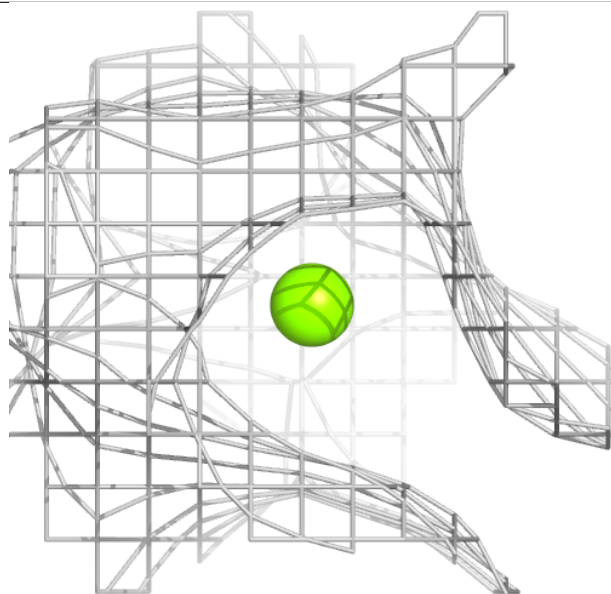
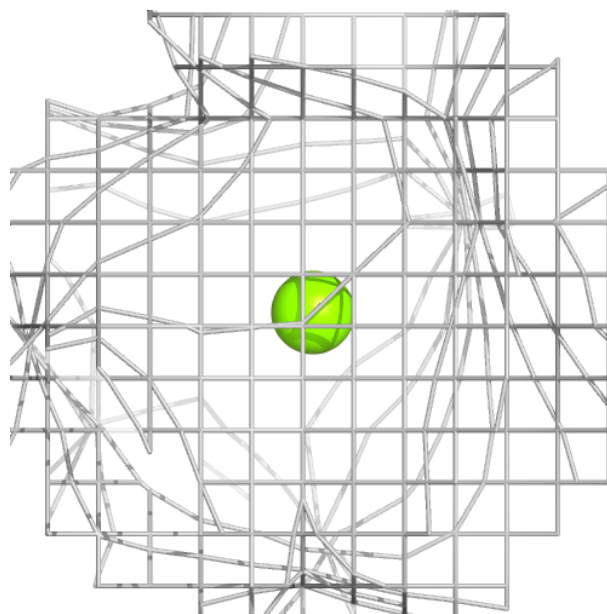
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q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

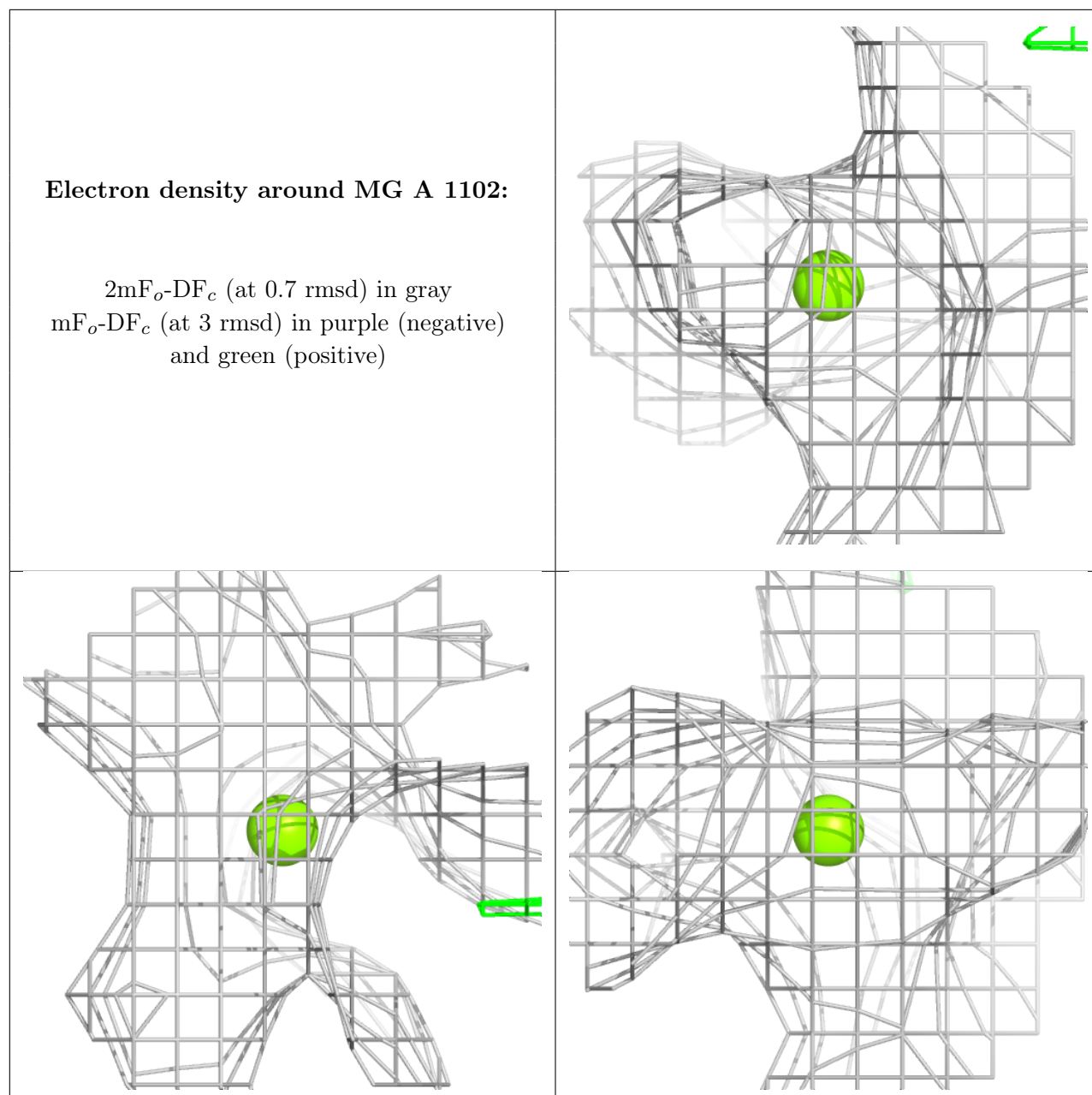
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
6	MG	A	1101	1/1	0.96	0.17	23,23,23,23	0
6	MG	A	1102	1/1	0.97	0.06	41,41,41,41	0
7	EDO	A	1103	4/4	0.97	0.21	38,42,46,49	0
6	MG	B	101	1/1	0.98	0.07	40,40,40,40	0

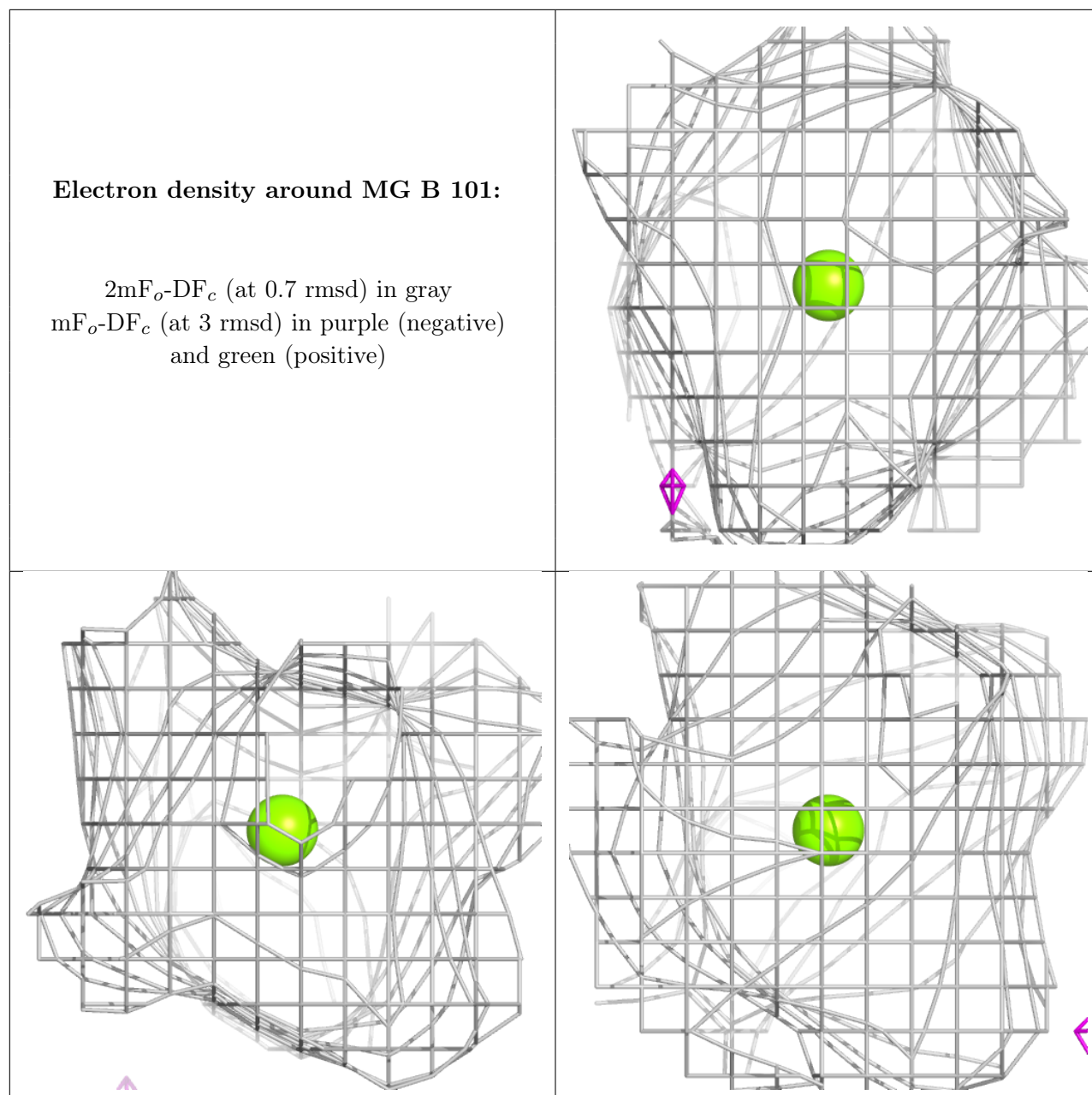
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

**Electron density around MG A 1101:**

$2mF_o-DF_c$  (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.