



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 20, 2026 – 05:50 PM UTC

PDB ID : 6LTU / pdb_00006ltu
Title : Crystal structure of Cas12i2 ternary complex with double Mg²⁺ 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

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>.

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

MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

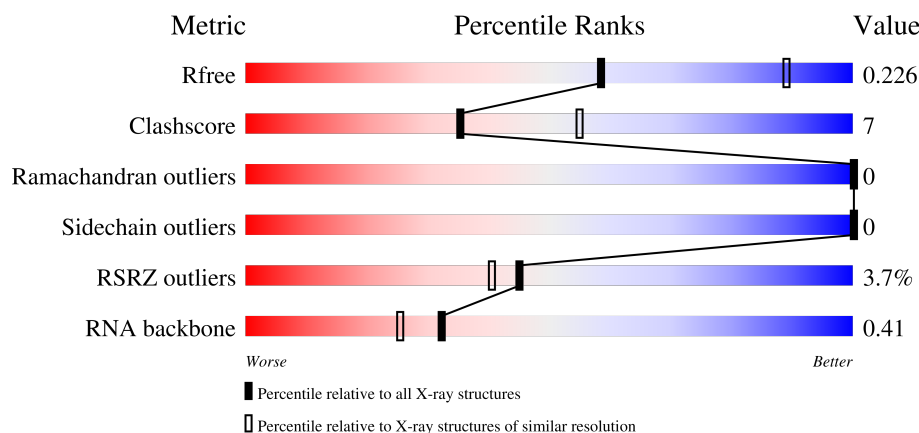
1 Overall quality at a glance

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}	180053	4770 (2.60-2.56)
Clashscore	190562	5124 (2.60-2.56)
Ramachandran outliers	187476	5046 (2.60-2.56)
Sidechain outliers	187428	5046 (2.60-2.56)
RSRZ outliers	180081	4770 (2.60-2.56)
RNA backbone	3983	1011 (2.82-2.34)

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 ≥ 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	4% 81% 17% .
2	B	58	45% 24% 14% . 16%
3	D	12	8% 42% 42% 8% 8%
4	C	35	6% 40% 57% .

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Mol	Chain	Length	Quality of chain
5	E	5	 80%20%

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
1	A	1034	Total	C	N	O	S	0	1	0
			8312	5280	1467	1538	27			

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

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

- 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
3	D	11	Total	C	N	O	P	0	0	0
			217	106	32	69	10			

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

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

- Molecule 5 is a DNA chain called trans ssDNA.

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

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

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

- Molecule 7 is 1,2-ETHANEDIOL (CCD ID: EDO) (formula: $C_2H_6O_2$).



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

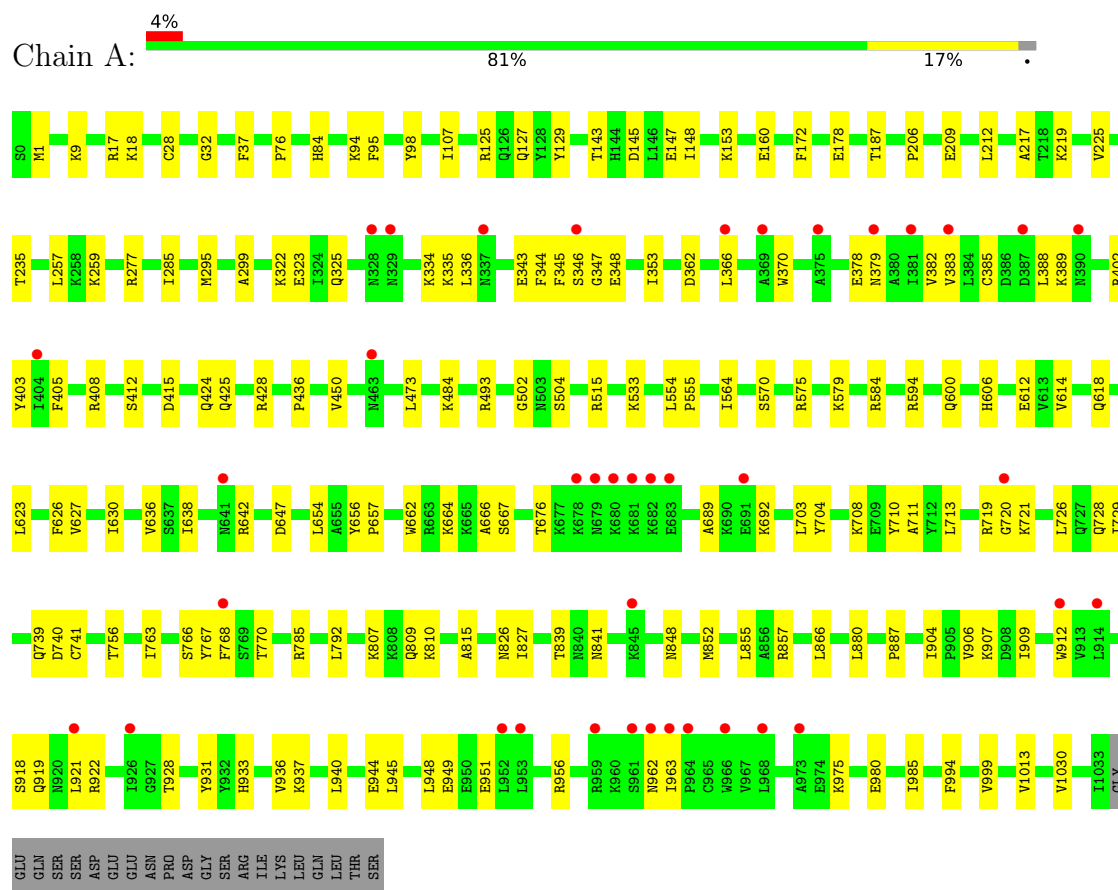
- Molecule 8 is water.

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

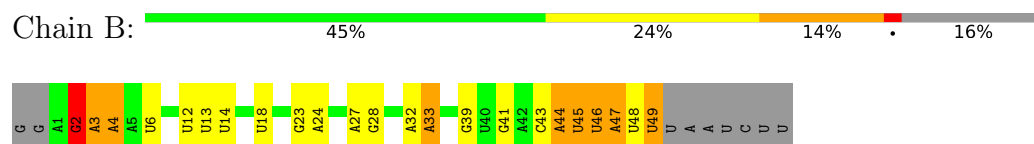
3 Residue-property plots [i](#)

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.

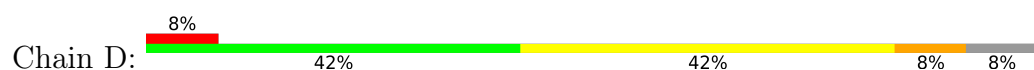
• Molecule 1: Cas12i2



• Molecule 2: RNA (56-mer)

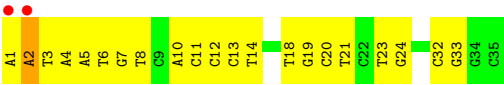


• Molecule 3: DNA (5'-D(*GP*CP*CP*GP*CP*TP*TP*TP*CP*TP*T)-3')

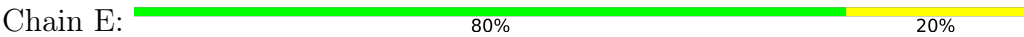




● Molecule 4: DNA (35-MER)



● Molecule 5: trans ssDNA



4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	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) 96.9 (46.87-2.57)	Depositor EDS
R_{merge}	0.16	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.76 (at 2.58Å)	Xtriage
Refinement program	PHENIX 1.18.2_3874	Depositor
R, R_{free}	0.202 , 0.229 (Not available) , 0.226	Depositor DCC
R_{free} test set	2525 reflections (4.73%)	wwPDB-VP
Wilson B-factor (Å ²)	41.7	Xtriage
Anisotropy	0.205	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 31.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	10704	wwPDB-VP
Average B, all atoms (Å ²)	49.0	wwPDB-VP

Xtriage'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.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: 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	0.49	7/8481 (0.1%)	0.61	5/11451 (0.0%)
2	B	0.31	0/1170	0.91	8/1822 (0.4%)
3	D	0.39	0/240	0.81	1/368 (0.3%)
4	C	0.28	0/800	0.86	6/1232 (0.5%)
5	E	0.46	0/91	0.88	0/138
All	All	0.46	7/10782 (0.1%)	0.68	20/15011 (0.1%)

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	579	LYS	C-O	-8.02	1.14	1.23
1	A	919	GLN	C-O	-7.26	1.15	1.24
1	A	618	GLN	C-O	-6.16	1.16	1.24
1	A	726	LEU	C-O	-5.91	1.17	1.24
1	A	383	VAL	C-N	5.89	1.42	1.33
1	A	739	GLN	C-O	-5.16	1.17	1.23
1	A	729	ILE	C-O	-5.13	1.17	1.24

All (20) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	3	A	C4'-C3'-O3'	-15.36	86.36	109.40
2	B	47	A	C1'-C2'-O2'	-15.00	85.90	108.40
1	A	719	ARG	N-CA-C	11.41	123.27	111.07
2	B	2	G	C4'-C3'-O3'	-10.37	93.85	109.40
4	C	2	DA	C4-N9-C1'	-9.99	112.06	127.05
4	C	2	DA	C8-N9-C1'	9.90	141.91	127.05
3	D	10	DT	C2'-C3'-O3'	-9.39	97.41	111.50
2	B	47	A	C4'-C3'-O3'	8.86	126.29	113.00
2	B	4	A	C1'-C2'-O2'	-8.80	95.21	108.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	344	PHE	N-CA-C	-8.74	100.54	114.09
4	C	2	DA	C4'-C3'-O3'	-7.70	98.45	110.00
1	A	721	LYS	N-CA-C	6.93	119.67	110.53
4	C	3	DT	C2'-C3'-O3'	-6.54	101.68	111.50
2	B	47	A	N9-C1'-C2'	-6.37	102.45	112.00
4	C	2	DA	C2'-C3'-O3'	6.26	120.89	111.50
2	B	4	A	C4'-C3'-O3'	6.12	122.18	113.00
2	B	47	A	C3'-C2'-C1'	5.95	107.25	101.30
4	C	2	DA	N9-C1'-C2'	5.70	122.05	113.50
1	A	689	ALA	N-CA-C	5.26	118.48	111.75
1	A	728	GLN	CA-C-O	5.24	126.10	120.55

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	121	0
2	B	1045	0	529	16	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	4	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	147	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 (147) 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:1:MET:HE2	1:A:584:ARG:HH11	1.46	0.80
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:1:MET:HE2	1:A:584:ARG:NH1	2.07	0.69
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
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:921:LEU:HD23	1:A:921:LEU:C	2.29	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:918:SER:C	1:A:922:ARG:NH1	2.64	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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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: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:907:LYS:NZ	1:A:980:GLU:OE2	2.37	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
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:921:LEU:C	1:A:921:LEU:CD2	2.88	0.47
1:A:810:LYS:NZ	8:A:1216:HOH:O	2.46	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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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:704:TYR:OH	2:B:12:U:OP1	2.31	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:336:LEU:C	1:A:336:LEU:HD23	2.42	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:335:LYS:O	1:A:403:TYR:OH	2.30	0.43
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
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
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:366:LEU:C	1:A:366:LEU:HD23	2.44	0.42
1:A:912:TRP:H	1:A:912:TRP:CD1	2.37	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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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
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	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	890/931 (96%)	890 (100%)	0	100	100

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 (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	19	ASN
1	A	90	ASN
1	A	163	GLN
1	A	329	ASN
1	A	459	ASN
1	A	620	HIS
1	A	641	ASN
1	A	826	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

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Mol	Chain	Res	Type
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 oligosaccharides 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	0.43	0	2,2,2	0.40	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, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	1034/1055 (98%)	0.32	39 (3%) 44 39	25, 45, 85, 107	1 (0%)
2	B	49/58 (84%)	-0.39	0 100 100	32, 40, 108, 141	0
3	D	11/12 (91%)	-0.02	1 (9%) 15 12	36, 39, 62, 97	0
4	C	35/35 (100%)	0.01	2 (5%) 29 25	35, 45, 129, 140	0
5	E	5/5 (100%)	0.08	0 100 100	46, 49, 59, 69	0
All	All	1134/1165 (97%)	0.28	42 (3%) 45 40	25, 45, 87, 141	1 (0%)

All (42) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	961	SER	4.8
1	A	679	ASN	4.1
1	A	966	TRP	4.0
3	D	11	DT	3.4
1	A	328	ASN	3.1
1	A	926	ILE	3.1
1	A	768	PHE	3.0
1	A	369	ALA	2.9
1	A	953	LEU	2.9
1	A	720	GLY	2.9
1	A	383	VAL	2.8
1	A	329	ASN	2.8
1	A	387	ASP	2.8
1	A	366	LEU	2.7
1	A	921	LEU	2.7
1	A	375	ALA	2.6
1	A	845	LYS	2.6
1	A	404	ILE	2.6
1	A	683	GLU	2.5
1	A	682	LYS	2.5

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Mol	Chain	Res	Type	RSRZ
1	A	912	TRP	2.5
1	A	959	ARG	2.5
1	A	968	LEU	2.4
4	C	2	DA	2.4
1	A	964	PRO	2.3
4	C	1	DA	2.3
1	A	952	LEU	2.3
1	A	691	GLU	2.3
1	A	379	ASN	2.2
1	A	337	ASN	2.2
1	A	962	ASN	2.2
1	A	681	LYS	2.2
1	A	381	ILE	2.1
1	A	346	SER	2.1
1	A	678	LYS	2.1
1	A	963	ILE	2.1
1	A	680	LYS	2.1
1	A	914	LEU	2.1
1	A	390	ASN	2.0
1	A	641	ASN	2.0
1	A	973	ALA	2.0
1	A	463	ASN	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 oligosaccharides in this entry.

6.4 Ligands [i](#)

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

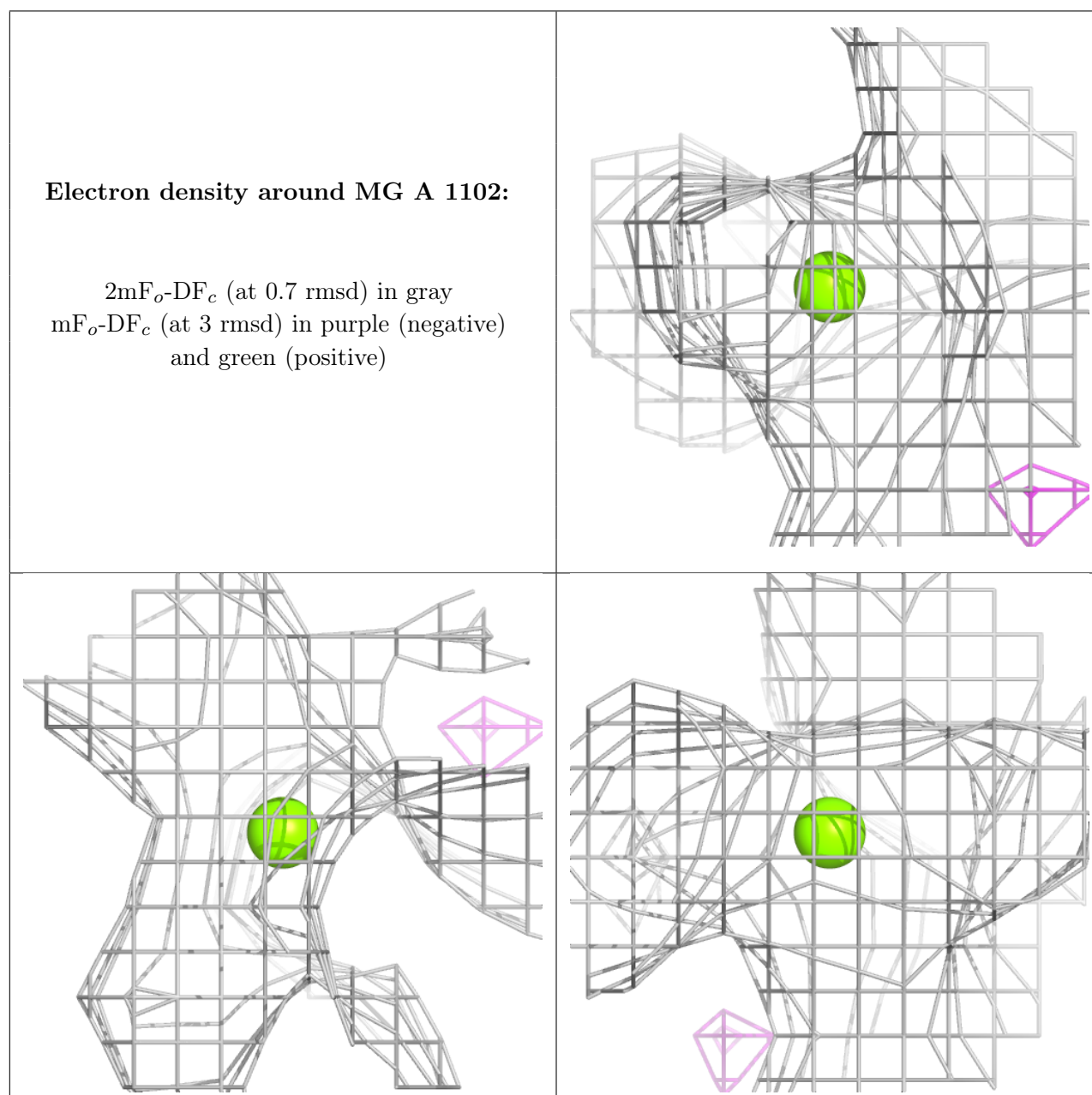
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
7	EDO	A	1103	4/4	0.93	0.10	38,42,46,49	0

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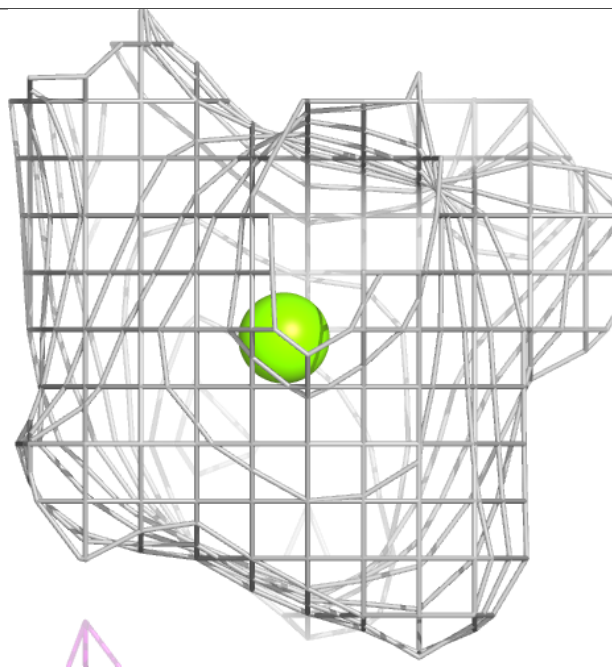
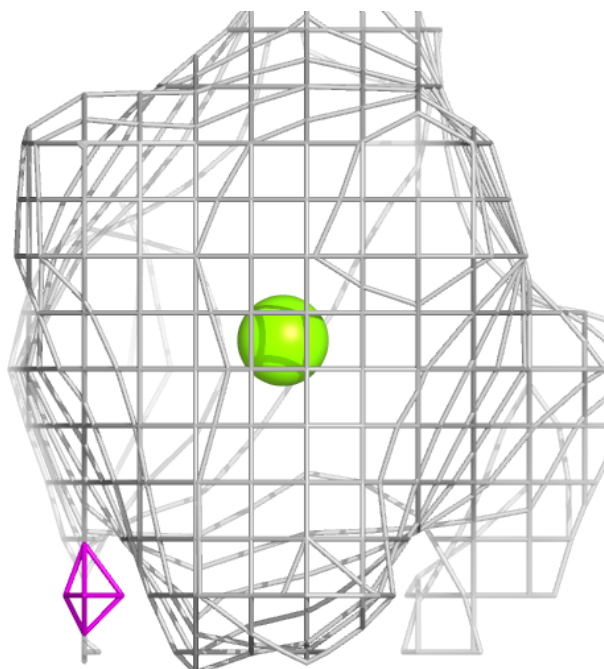
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
6	MG	A	1102	1/1	0.97	0.09	41,41,41,41	0
6	MG	B	101	1/1	0.99	0.03	40,40,40,40	0
6	MG	A	1101	1/1	0.99	0.05	23,23,23,23	0

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.



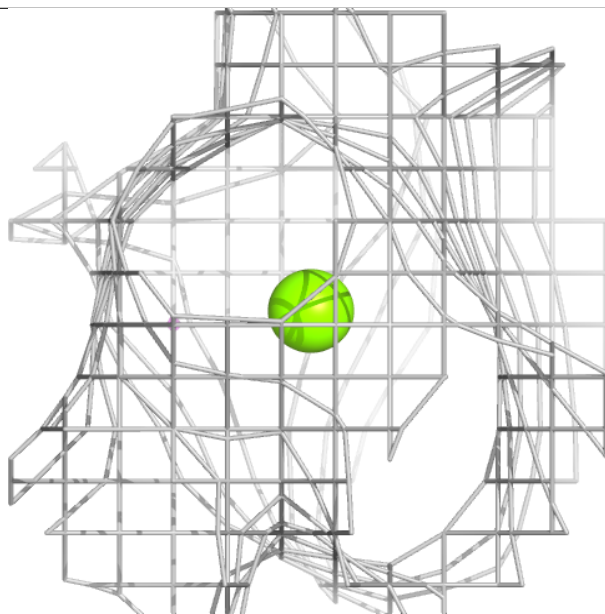
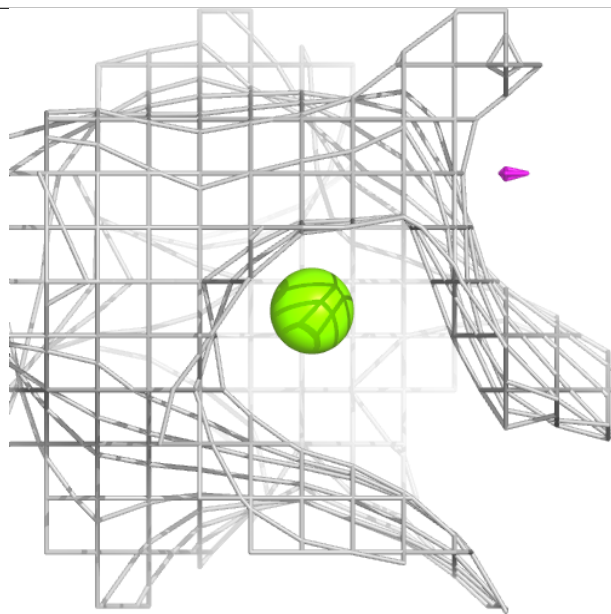
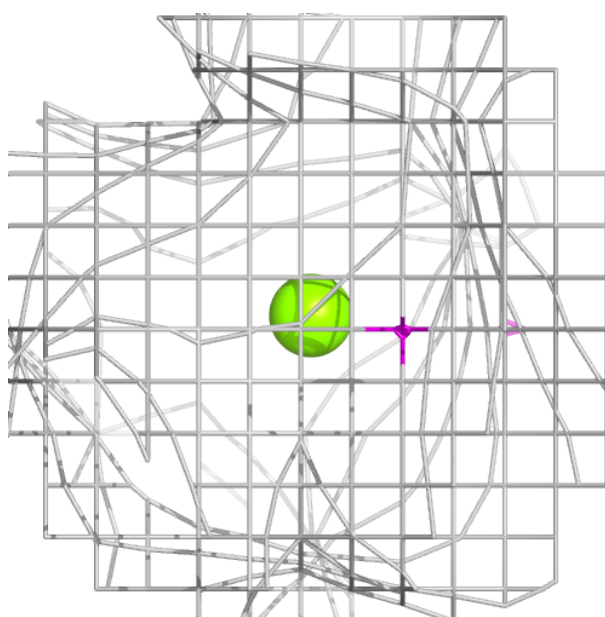
Electron density around MG B 101:

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



Electron density around 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 ⓘ

There are no such residues in this entry.