



wwPDB X-ray Structure Validation Summary Report ⓘ

Aug 22, 2020 – 09:24 AM BST

PDB ID : 6PXU
Title : Crystal structure of human GalNAc-T12 bound to a diglycosylated peptide, Mn²⁺, and UDP
Authors : Samara, N.L.; Fernandez, A.J.
Deposited on : 2019-07-27
Resolution : 2.01 Å(reported)

This is a wwPDB X-ray Structure Validation Summary 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 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 : 2.13.1
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) : 2.13.1

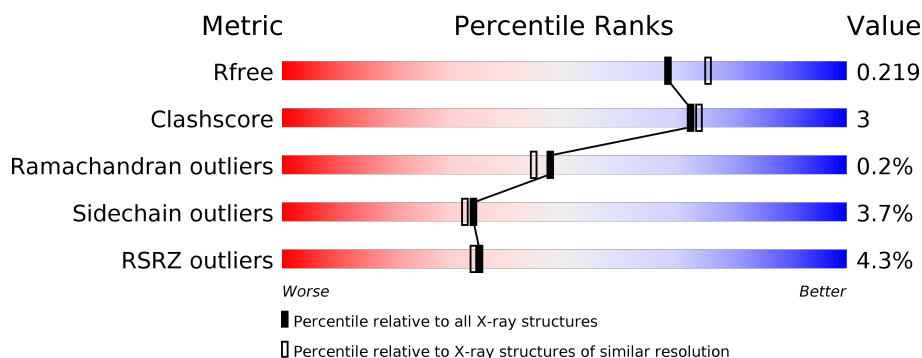
1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.01 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	8085 (2.00-2.00)
Clashscore	141614	9178 (2.00-2.00)
Ramachandran outliers	138981	9054 (2.00-2.00)
Sidechain outliers	138945	9053 (2.00-2.00)
RSRZ outliers	127900	7900 (2.00-2.00)

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 ≥ 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	543	<div> <div>4%</div> <div> <div></div> <div>87%</div> <div>10%</div> <div>•</div> </div> </div>
1	B	543	<div> <div>3%</div> <div> <div></div> <div>88%</div> <div>9%</div> <div>••</div> </div> </div>
2	C	21	<div> <div>29%</div> <div> <div>52%</div> <div>10%</div> <div>38%</div> </div> </div>
2	D	21	<div> <div>29%</div> <div> <div>62%</div> <div>5%</div> <div>5%</div> <div>29%</div> </div> </div>

2 Entry composition [i](#)

There are 8 unique types of molecules in this entry. The entry contains 9719 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.

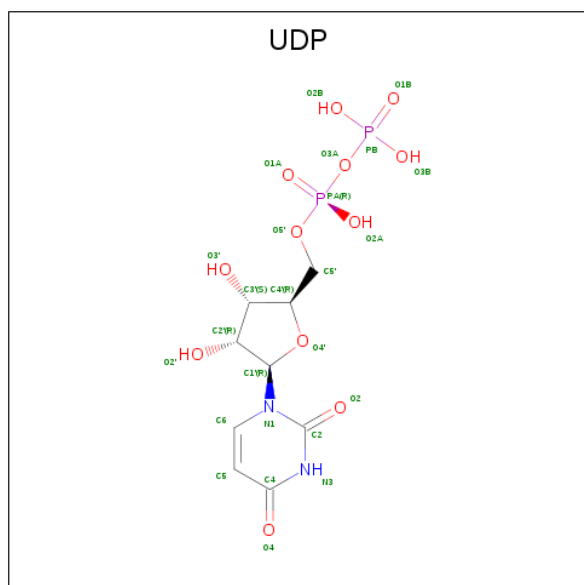
- Molecule 1 is a protein called Polypeptide N-acetylgalactosaminyltransferase 12.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	526	Total	C	N	O	S	0	10	0
			4351	2761	770	793	27			
1	B	532	Total	C	N	O	S	0	8	0
			4405	2793	787	801	24			

- Molecule 2 is a protein called GAGATGAGAGYYITPRTGAGA.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	C	13	Total	C	N	O	0	0	0
			93	59	16	18			
2	D	15	Total	C	N	O	0	0	0
			102	64	18	20			

- Molecule 3 is URIDINE-5'-DIPHOSPHATE (three-letter code: UDP) (formula: $C_9H_{14}N_2O_{12}P_2$) (labeled as "Ligand of Interest" by author).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	N	O	P	0	0
			25	9	2	12	2		
3	B	1	Total	C	N	O	P	0	0
			25	9	2	12	2		

- Molecule 4 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula: C₂H₆O₂).



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

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

- Molecule 5 is GLYCEROL (three-letter code: GOL) (formula: $C_3H_8O_3$).

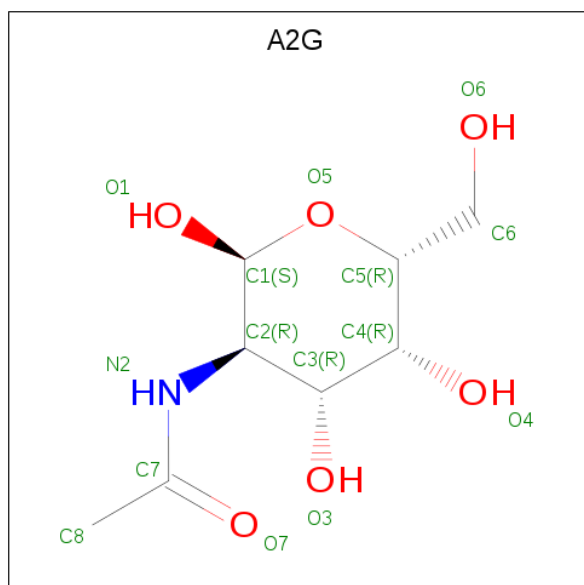


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	C	O	0	0
			6	3	3		
5	A	1	Total	C	O	0	0
			6	3	3		
5	A	1	Total	C	O	0	0
			6	3	3		
5	B	1	Total	C	O	0	0
			6	3	3		
5	B	1	Total	C	O	0	0
			6	3	3		
5	B	1	Total	C	O	0	0
			6	3	3		
5	B	1	Total	C	O	0	0
			6	3	3		

- Molecule 6 is MANGANESE (II) ION (three-letter code: MN) (formula: Mn) (labeled as "Ligand of Interest" by author).

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

- Molecule 7 is 2-acetamido-2-deoxy- α -D-galactopyranose (three-letter code: A2G) (formula: $C_8H_{15}NO_6$) (labeled as "Ligand of Interest" by author).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
7	C	1	Total C N O 14 8 1 5	0	0
7	C	1	Total C N O 14 8 1 5	0	0
7	D	1	Total C N O 14 8 1 5	0	0
7	D	1	Total C N O 14 8 1 5	0	0

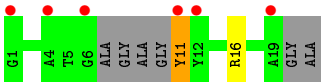
- Molecule 8 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
8	A	271	Total O 271 271	0	0
8	B	280	Total O 280 280	0	0
8	C	6	Total O 6 6	0	0

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	D	5	Total	O	0	0
			5	5		



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	72.82Å 73.12Å 74.44Å 113.08° 100.50° 108.23°	Depositor
Resolution (Å)	20.03 – 2.01 20.03 – 2.01	Depositor EDS
% Data completeness (in resolution range)	97.1 (20.03-2.01) 97.1 (20.03-2.01)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.67 (at 2.01Å)	Xtriage
Refinement program	PHENIX 1.10.1 _2155	Depositor
R, R_{free}	0.171 , 0.219 0.172 , 0.219	Depositor DCC
R_{free} test set	3944 reflections (4.83%)	wwPDB-VP
Wilson B-factor (Å ²)	29.3	Xtriage
Anisotropy	0.542	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 49.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	9719	wwPDB-VP
Average B, all atoms (Å ²)	42.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.07% 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: GOL, UDP, MN, EDO, A2G

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	0.40	0/4492	0.58	1/6085 (0.0%)
1	B	0.40	0/4545	0.58	2/6153 (0.0%)
2	C	0.33	0/94	0.52	0/126
2	D	0.35	0/103	0.66	0/138
All	All	0.40	0/9234	0.58	3/12502 (0.0%)

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	212	ARG	NE-CZ-NH2	-5.82	117.39	120.30
1	A	212	ARG	NE-CZ-NH1	-5.46	117.57	120.30
1	B	520	LEU	CA-CB-CG	5.24	127.34	115.30

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	4351	0	4238	28	0
1	B	4405	0	4307	31	0
2	C	93	0	85	1	0
2	D	102	0	96	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	A	25	0	11	3	0
3	B	25	0	11	0	0
4	A	20	0	30	1	0
4	B	28	0	42	5	0
5	A	18	0	24	0	0
5	B	30	0	40	3	0
6	A	2	0	0	0	0
6	B	2	0	0	0	0
7	C	28	0	24	1	0
7	D	28	0	24	0	0
8	A	271	0	0	3	0
8	B	280	0	0	4	0
8	C	6	0	0	0	0
8	D	5	0	0	0	0
All	All	9719	0	8932	62	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.

The worst 5 of 62 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:186:ARG:HH22	4:B:611:EDO:H11	1.57	0.69
1:A:304:VAL:HG22	1:A:354:GLU:HB3	1.77	0.65
1:A:208:LEU:HD22	3:A:601:UDP:H5'1	1.79	0.63
1:B:59:GLY:O	1:B:61[A]:ARG:NH2	2.33	0.61
1:B:131:ASP:HB3	4:B:613:EDO:H21	1.86	0.58

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	531/543 (98%)	518 (98%)	12 (2%)	1 (0%)	47	44
1	B	538/543 (99%)	525 (98%)	12 (2%)	1 (0%)	47	44
2	C	9/21 (43%)	9 (100%)	0	0	100	100
2	D	11/21 (52%)	11 (100%)	0	0	100	100
All	All	1089/1128 (96%)	1063 (98%)	24 (2%)	2 (0%)	47	44

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	334	VAL
1	B	334	VAL

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	476/476 (100%)	459 (96%)	17 (4%)	35	34
1	B	481/476 (101%)	463 (96%)	18 (4%)	34	32
2	C	8/8 (100%)	7 (88%)	1 (12%)	4	2
2	D	8/8 (100%)	6 (75%)	2 (25%)	0	0
All	All	973/968 (100%)	935 (96%)	38 (4%)	34	30

5 of 38 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	85	GLN
1	B	230	HIS
2	C	11	TYR
1	B	93	LEU
1	B	256	VAL

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	189	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 30 ligands modelled in this entry, 4 are monoatomic - leaving 26 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	$\# Z > 2$
4	EDO	B	603	-	3,3,3	0.44	0	2,2,2	0.42	0
4	EDO	A	604	-	3,3,3	0.51	0	2,2,2	0.32	0
5	GOL	B	610	-	5,5,5	0.38	0	5,5,5	0.21	0
5	GOL	A	607	-	5,5,5	0.35	0	5,5,5	0.46	0
5	GOL	B	607	-	5,5,5	0.38	0	5,5,5	0.35	0
4	EDO	B	605	-	3,3,3	0.50	0	2,2,2	0.26	0
7	A2G	C	101	2	14,14,15	0.45	0	17,19,21	0.81	0
4	EDO	B	612	-	3,3,3	0.47	0	2,2,2	0.30	0
7	A2G	D	102	2	14,14,15	0.48	0	17,19,21	2.28	4 (23%)
4	EDO	B	604	-	3,3,3	0.43	0	2,2,2	0.43	0
7	A2G	D	101	2	14,14,15	0.45	0	17,19,21	0.93	1 (5%)
3	UDP	A	601	6	20,26,26	1.32	2 (10%)	25,40,40	1.48	4 (16%)
5	GOL	B	609	-	5,5,5	0.37	0	5,5,5	0.20	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	UDP	B	602	6	20,26,26	1.11	1 (5%)	25,40,40	0.83	0
5	GOL	A	605	-	5,5,5	0.37	0	5,5,5	0.19	0
5	GOL	B	606	-	5,5,5	0.42	0	5,5,5	0.71	0
5	GOL	B	608	-	5,5,5	0.43	0	5,5,5	0.65	0
4	EDO	B	613	-	3,3,3	0.44	0	2,2,2	0.45	0
5	GOL	A	606	-	5,5,5	0.43	0	5,5,5	0.33	0
4	EDO	A	609	-	3,3,3	0.46	0	2,2,2	0.37	0
4	EDO	A	602	-	3,3,3	0.52	0	2,2,2	0.22	0
4	EDO	A	608	-	3,3,3	0.49	0	2,2,2	0.37	0
4	EDO	B	611	-	3,3,3	0.50	0	2,2,2	0.26	0
4	EDO	B	614	-	3,3,3	0.46	0	2,2,2	0.31	0
7	A2G	C	102	2	14,14,15	0.51	0	17,19,21	0.83	0
4	EDO	A	603	-	3,3,3	0.52	0	2,2,2	0.26	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
4	EDO	B	603	-	-	0/1/1/1	-
4	EDO	A	604	-	-	0/1/1/1	-
5	GOL	B	610	-	-	2/4/4/4	-
5	GOL	A	607	-	-	0/4/4/4	-
5	GOL	B	607	-	-	0/4/4/4	-
4	EDO	B	605	-	-	0/1/1/1	-
7	A2G	C	101	2	-	0/6/23/26	0/1/1/1
4	EDO	B	612	-	-	0/1/1/1	-
7	A2G	D	102	2	-	5/6/23/26	0/1/1/1
4	EDO	B	604	-	-	0/1/1/1	-
7	A2G	D	101	2	-	2/6/23/26	0/1/1/1
3	UDP	A	601	6	-	7/14/32/32	0/2/2/2
5	GOL	B	609	-	-	2/4/4/4	-
3	UDP	B	602	6	-	3/14/32/32	0/2/2/2
5	GOL	A	605	-	-	4/4/4/4	-
5	GOL	B	606	-	-	2/4/4/4	-
5	GOL	B	608	-	-	1/4/4/4	-
4	EDO	B	613	-	-	0/1/1/1	-
5	GOL	A	606	-	-	0/4/4/4	-
4	EDO	A	609	-	-	1/1/1/1	-
4	EDO	A	602	-	-	0/1/1/1	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	EDO	A	608	-	-	0/1/1/1	-
4	EDO	B	611	-	-	1/1/1/1	-
4	EDO	B	614	-	-	0/1/1/1	-
7	A2G	C	102	2	-	2/6/23/26	0/1/1/1
4	EDO	A	603	-	-	1/1/1/1	-

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	601	UDP	C4-N3	3.19	1.38	1.33
3	B	602	UDP	C4-N3	3.05	1.38	1.33
3	A	601	UDP	O4'-C4'	-2.55	1.39	1.45

The worst 5 of 9 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
7	D	102	A2G	C4-C3-C2	-6.36	101.69	111.02
7	D	102	A2G	C2-N2-C7	4.26	128.97	122.90
3	A	601	UDP	O5'-PA-O1A	3.94	124.46	109.07
3	A	601	UDP	O4'-C4'-C5'	-3.38	98.26	109.37
7	D	102	A2G	O5-C1-C2	-3.19	106.25	111.29

There are no chirality outliers.

5 of 33 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	B	610	GOL	O1-C1-C2-O2
5	B	610	GOL	O1-C1-C2-C3
7	D	102	A2G	C3-C2-N2-C7
3	A	601	UDP	C2'-C1'-N1-C6
3	A	601	UDP	O4'-C1'-N1-C6

There are no ring outliers.

8 monomers are involved in 13 short contacts:

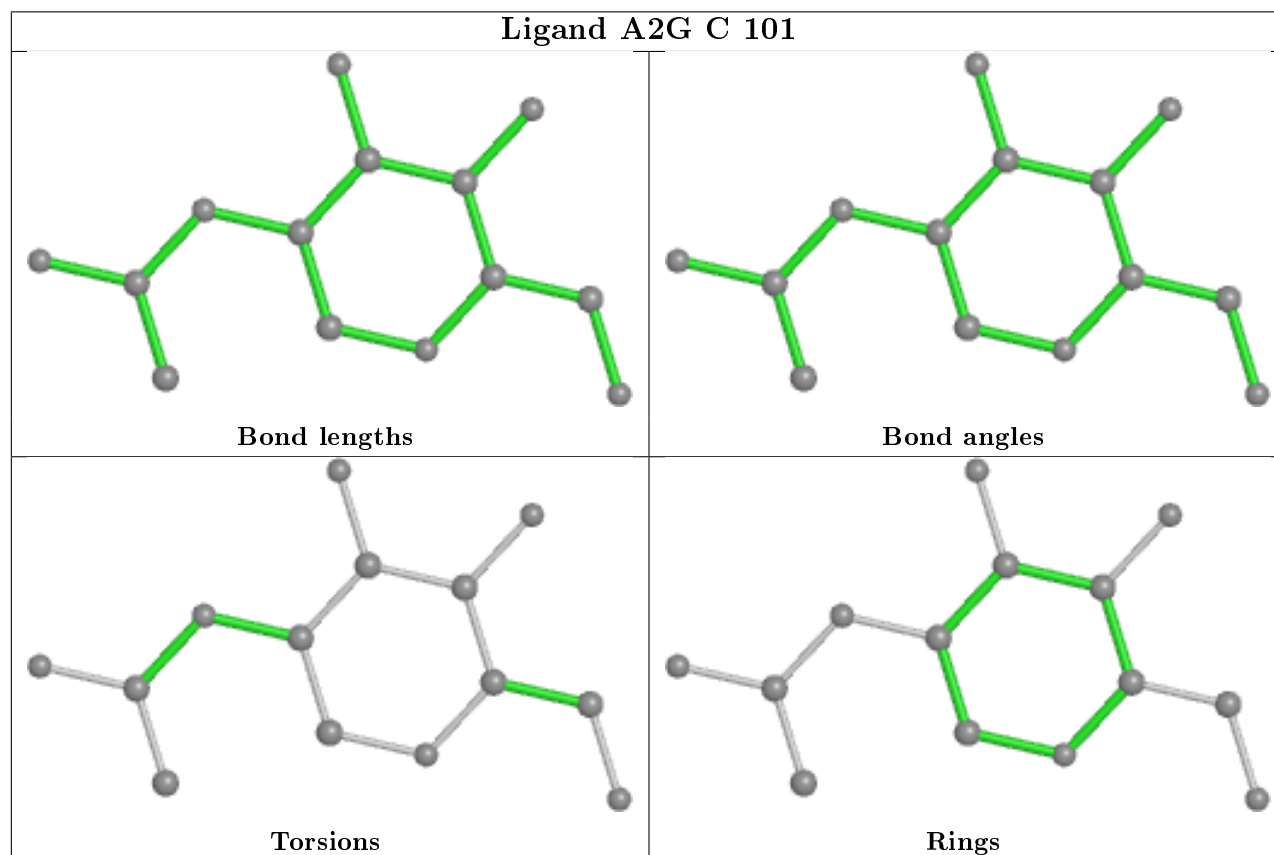
Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	B	610	GOL	2	0
3	A	601	UDP	3	0
5	B	606	GOL	1	0
4	B	613	EDO	2	0
4	A	608	EDO	1	0

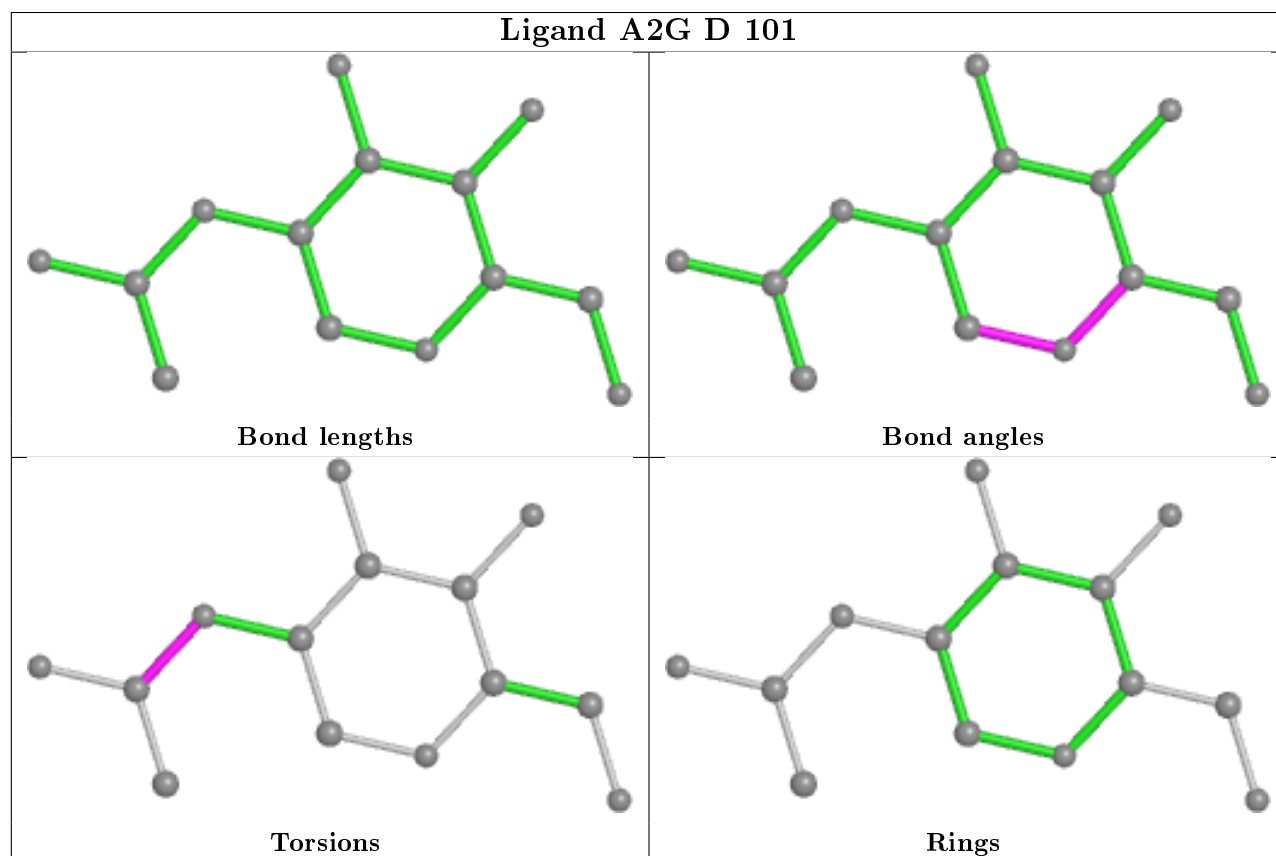
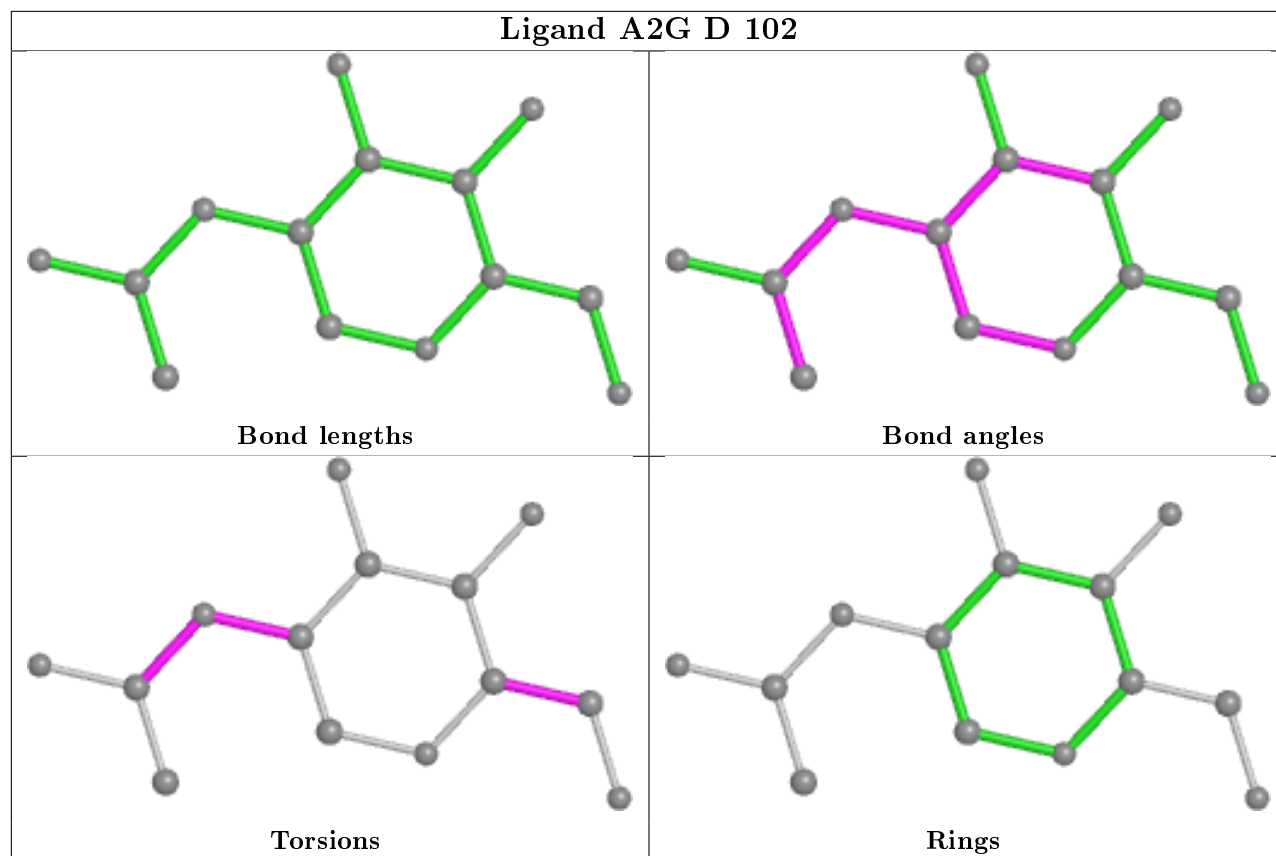
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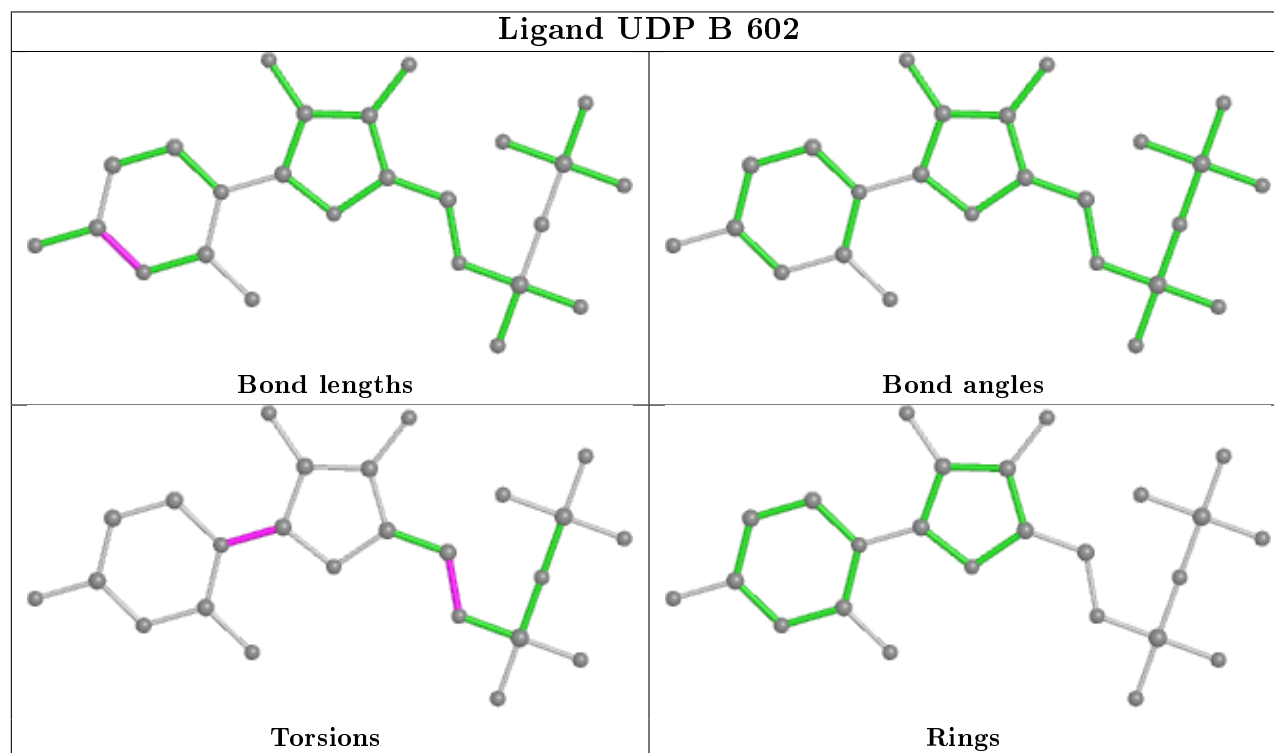
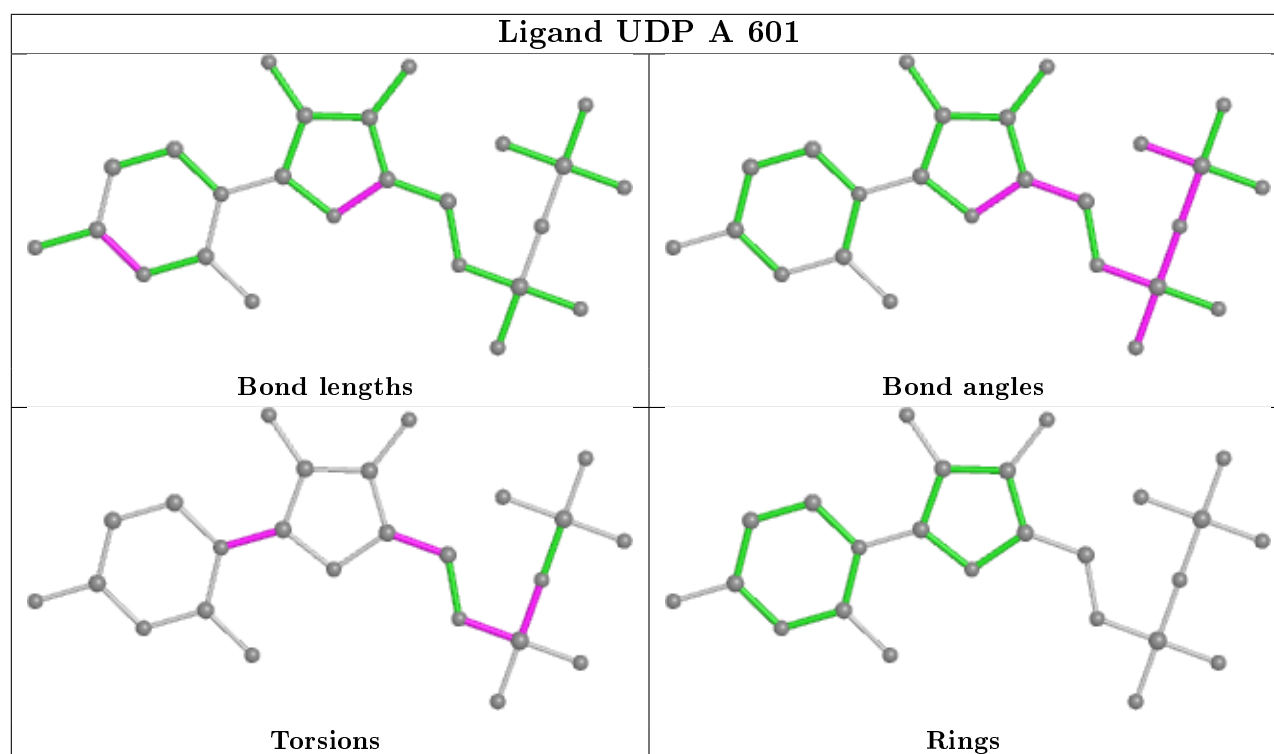
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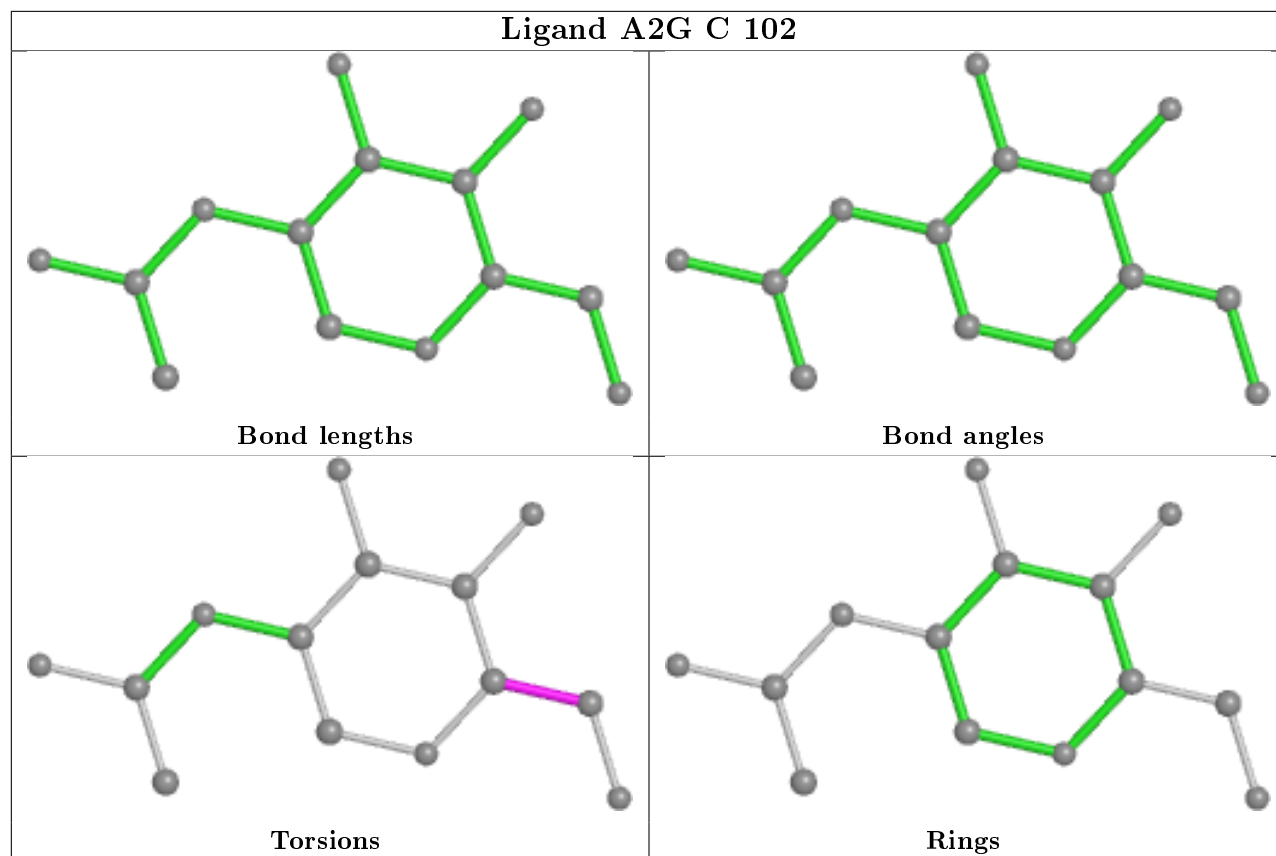
Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	B	611	EDO	1	0
4	B	614	EDO	2	0
7	C	102	A2G	1	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.









5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data [i](#)

6.1 Protein, DNA and RNA chains [i](#)

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 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	526/543 (96%)	-0.23	20 (3%)	40 39	20, 37, 75, 112	2 (0%)
1	B	532/543 (97%)	-0.29	15 (2%)	53 51	19, 36, 72, 112	0
2	C	13/21 (61%)	1.52	6 (46%)	0 0	39, 54, 96, 108	0
2	D	15/21 (71%)	1.52	6 (40%)	0 0	58, 68, 87, 96	0
All	All	1086/1128 (96%)	-0.21	47 (4%)	35 34	19, 37, 75, 112	2 (0%)

The worst 5 of 47 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	72	ALA	6.7
1	B	70	VAL	5.4
1	B	464	PRO	5.1
1	B	511	ALA	5.0
1	B	73	ASN	4.9

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](#)

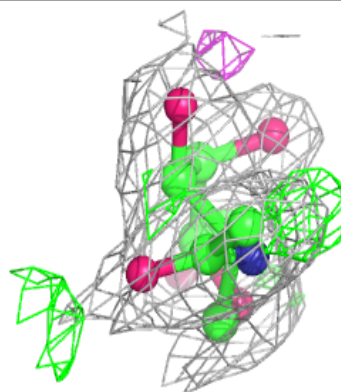
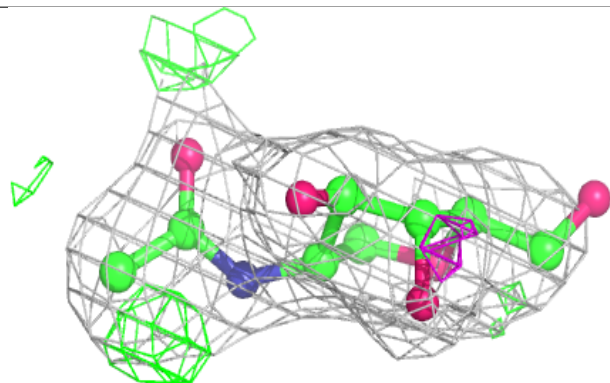
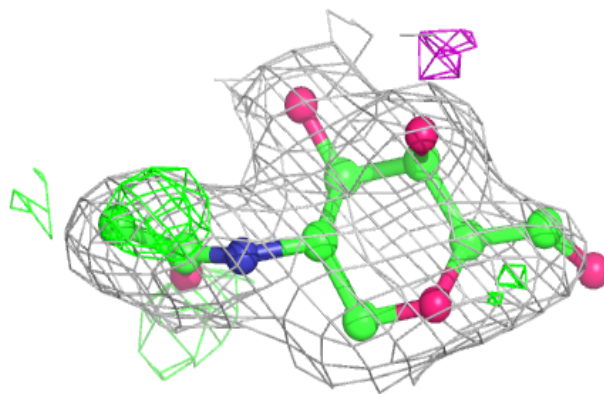
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled ‘Q< 0.9’ lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
4	EDO	A	603	4/4	0.68	0.31	64,66,71,72	0
4	EDO	B	611	4/4	0.82	0.26	61,66,71,73	0
5	GOL	B	609	6/6	0.83	0.19	60,72,74,74	0
5	GOL	A	607	6/6	0.84	0.23	49,63,67,71	0
4	EDO	A	602	4/4	0.85	0.16	46,54,59,61	0
5	GOL	B	610	6/6	0.85	0.18	55,65,70,70	0
4	EDO	A	609	4/4	0.85	0.18	76,77,77,79	0
4	EDO	B	613	4/4	0.87	0.46	62,63,64,66	0
5	GOL	B	608	6/6	0.88	0.15	35,39,48,52	0
5	GOL	B	607	6/6	0.88	0.23	58,63,64,65	0
4	EDO	B	603	4/4	0.88	0.15	67,68,70,74	0
7	A2G	D	102	14/15	0.89	0.14	28,51,64,72	0
4	EDO	B	612	4/4	0.89	0.24	68,68,69,69	0
4	EDO	B	605	4/4	0.90	0.23	65,66,68,68	0
4	EDO	A	608	4/4	0.91	0.16	58,60,61,62	0
4	EDO	A	604	4/4	0.91	0.11	46,46,50,56	0
4	EDO	B	614	4/4	0.91	0.25	52,52,53,56	0
5	GOL	A	606	6/6	0.91	0.12	40,47,48,49	0
5	GOL	A	605	6/6	0.92	0.14	49,58,62,62	0
7	A2G	C	102	14/15	0.93	0.15	45,63,71,73	0
6	MN	B	615	1/1	0.94	0.41	82,82,82,82	1
7	A2G	C	101	14/15	0.95	0.07	24,28,39,44	0
4	EDO	B	604	4/4	0.95	0.16	54,56,57,63	0
7	A2G	D	101	14/15	0.96	0.07	26,30,35,46	0
5	GOL	B	606	6/6	0.96	0.08	32,43,46,46	0
3	UDP	A	601	25/25	0.98	0.07	22,29,34,41	0
3	UDP	B	602	25/25	0.98	0.06	22,29,33,39	0
6	MN	A	611	1/1	0.99	0.06	31,31,31,31	1
6	MN	A	610	1/1	0.99	0.04	32,32,32,32	0
6	MN	B	601	1/1	1.00	0.02	30,30,30,30	0

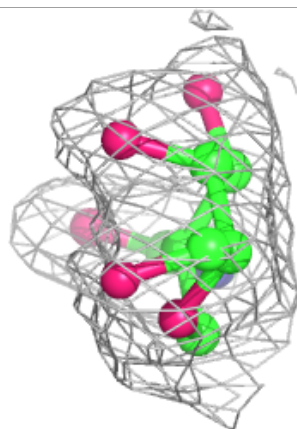
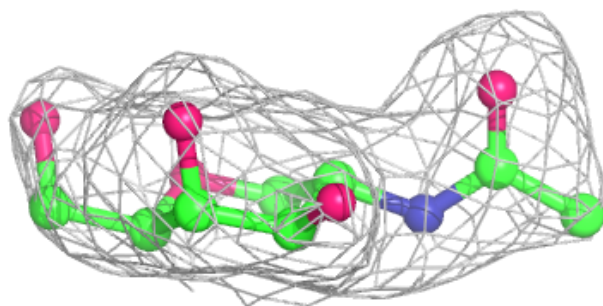
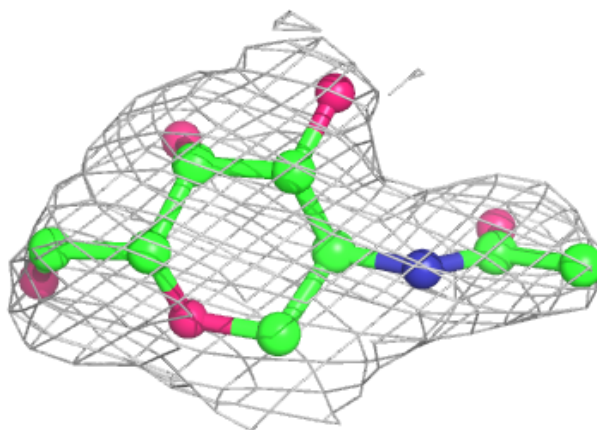
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around A2G D 102:

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

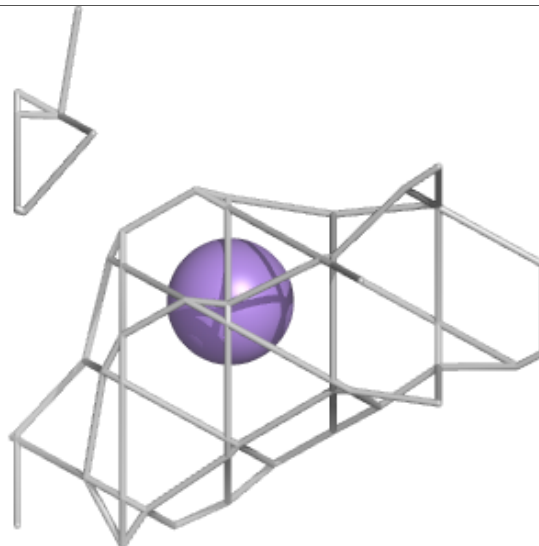
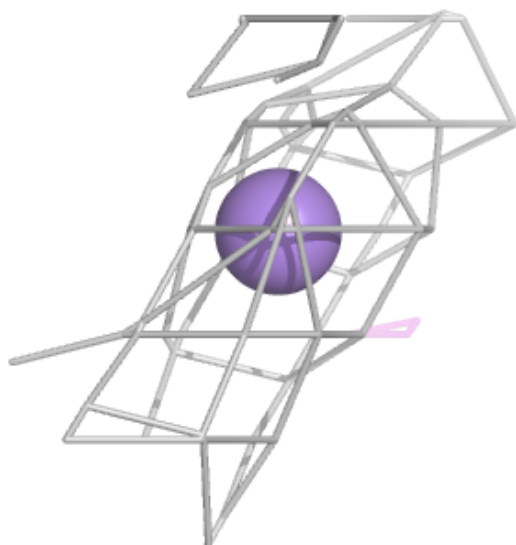
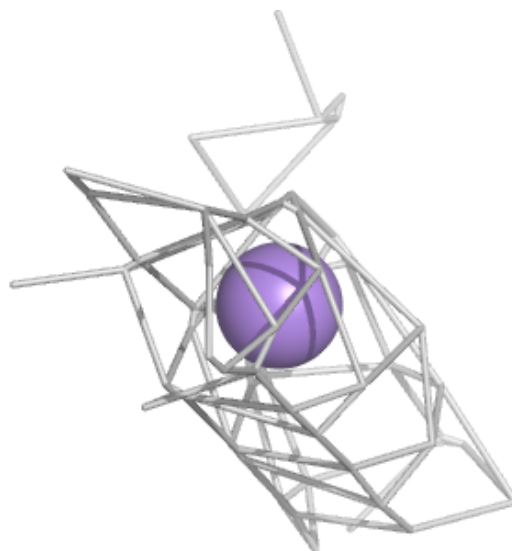
**Electron density around A2G C 102:**

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



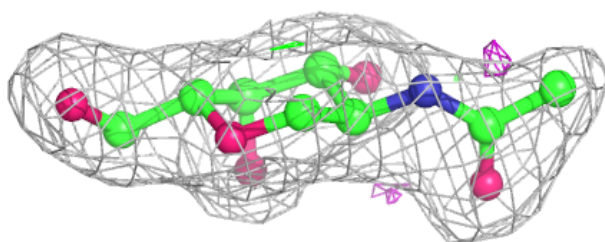
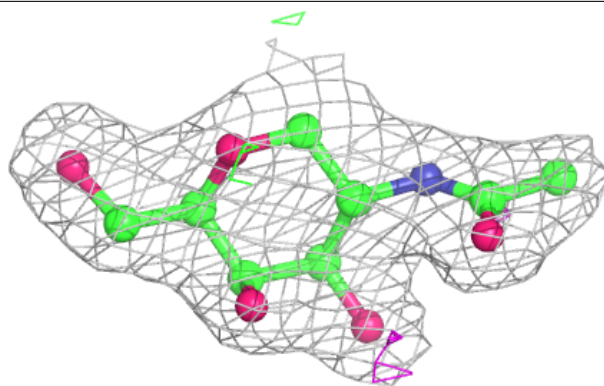
Electron density around MN B 615:

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

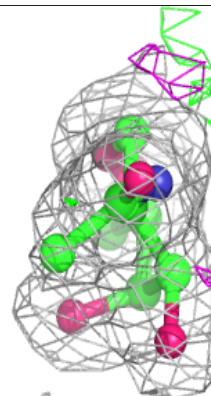
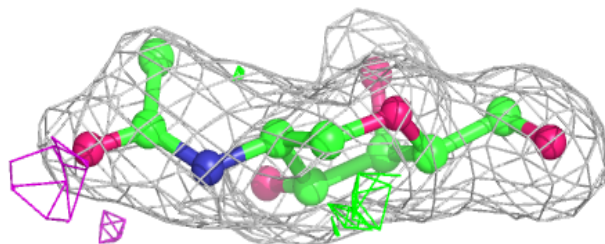
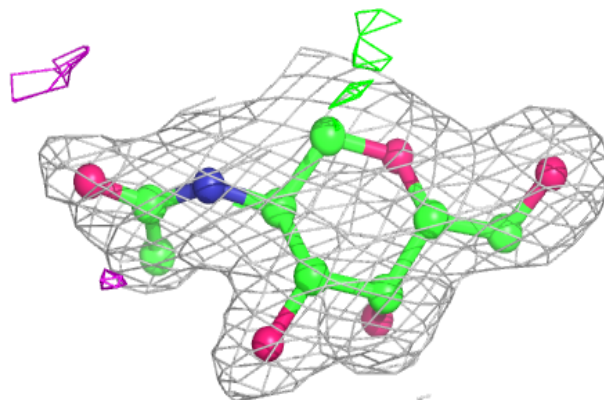


Electron density around A2G C 101:

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

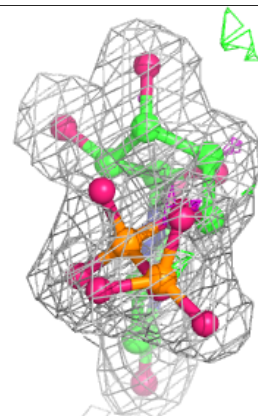
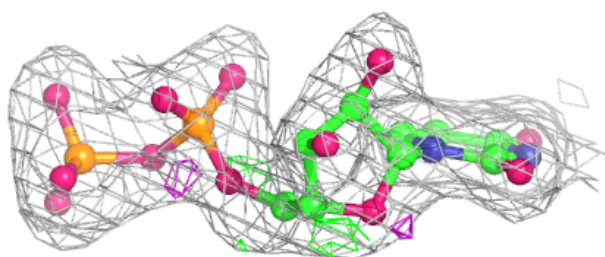
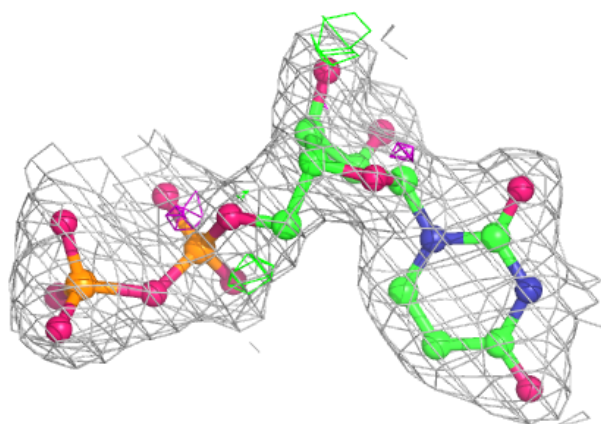
**Electron density around A2G D 101:**

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 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

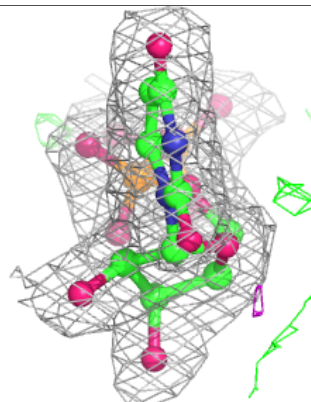
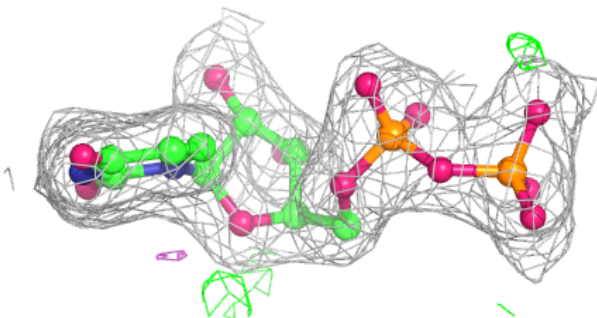
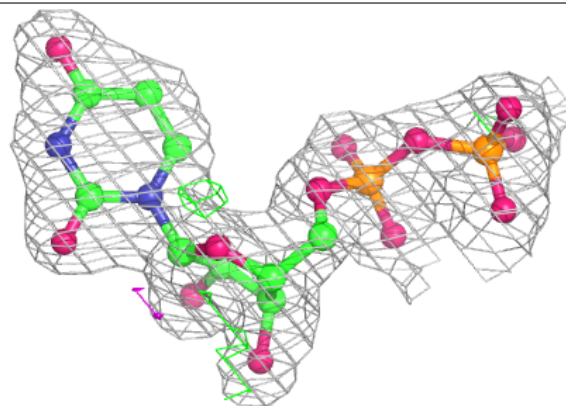


Electron density around UDP A 601:

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

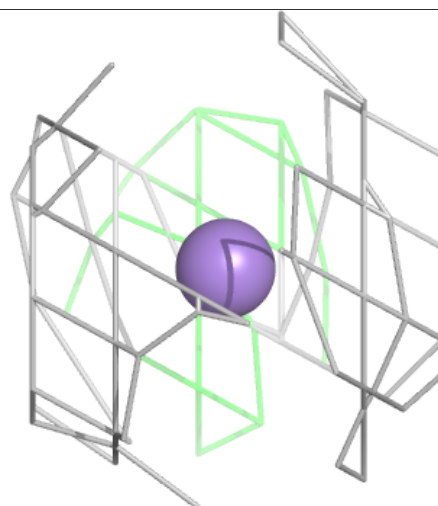
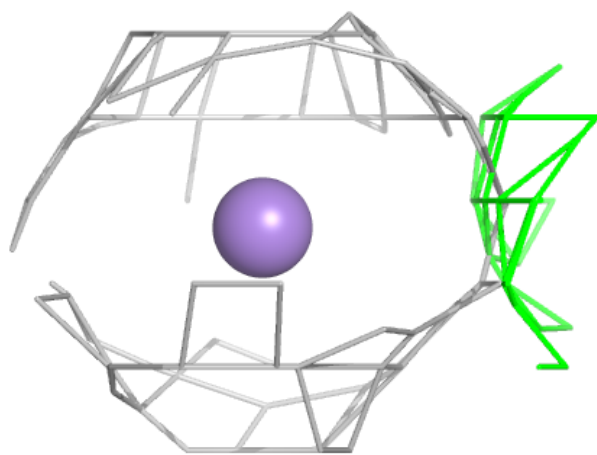
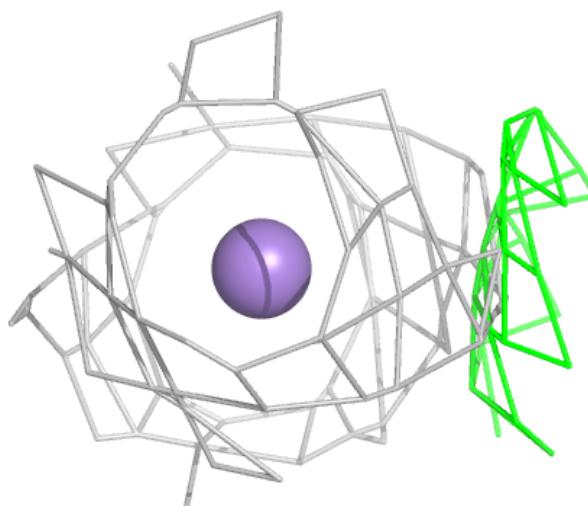
**Electron density around UDP B 602:**

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



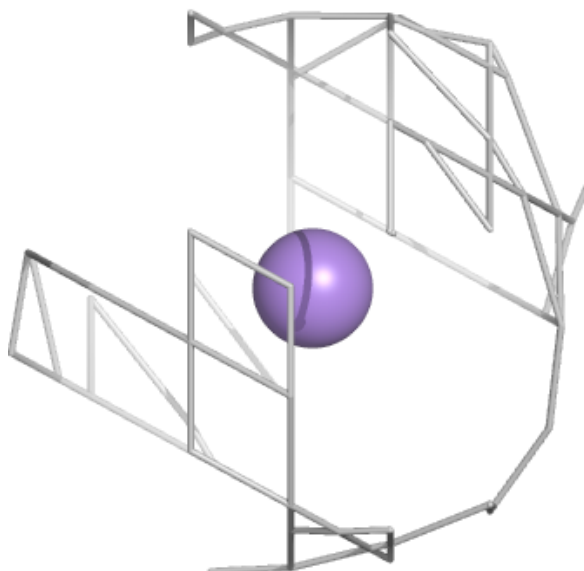
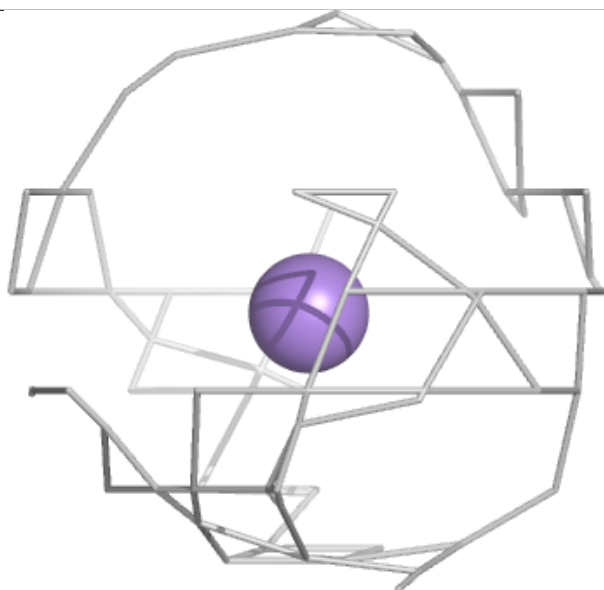
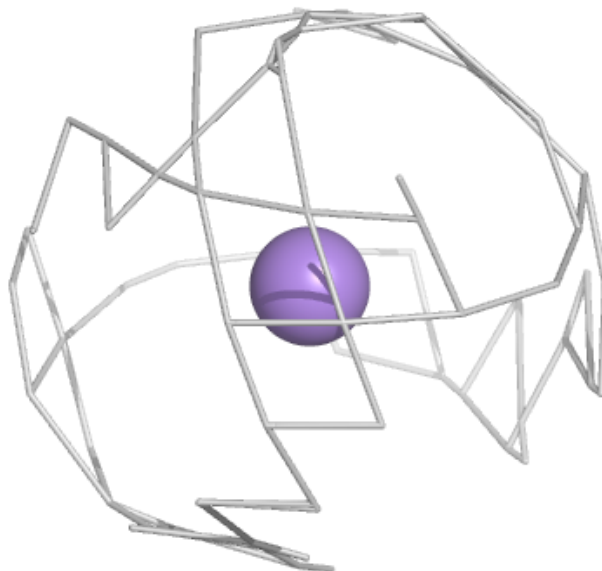
Electron density around MN A 611:

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



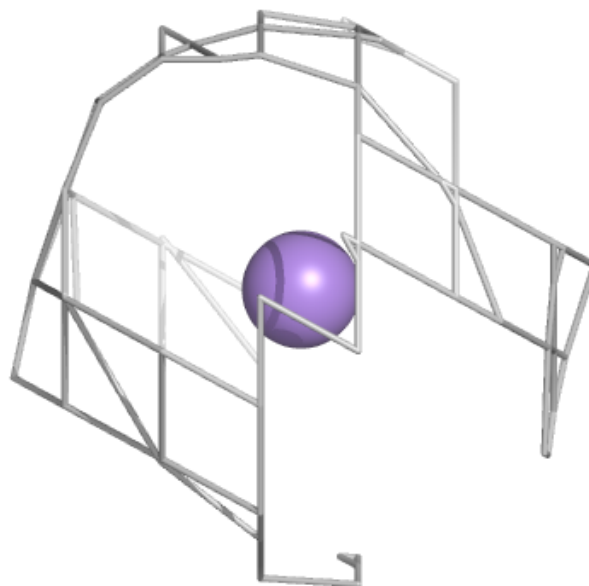
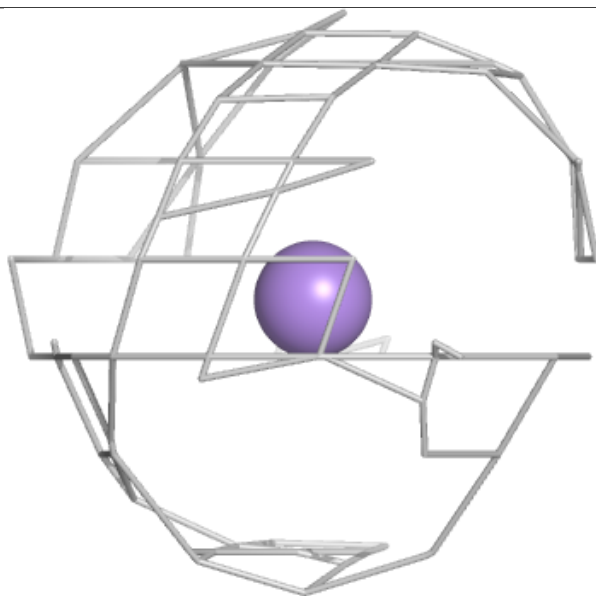
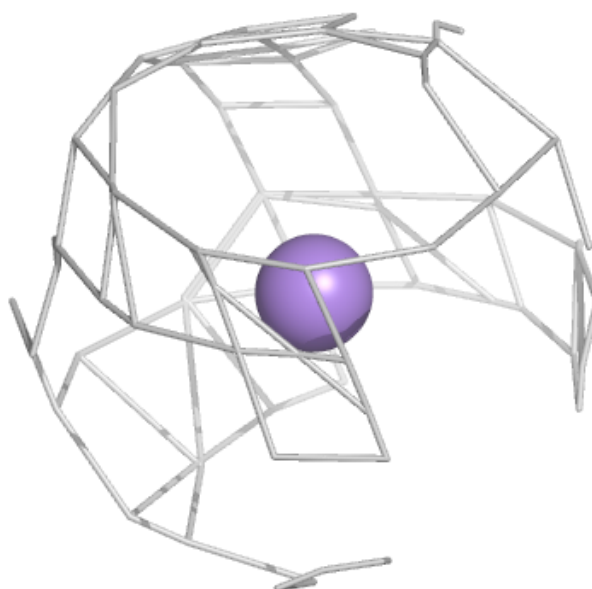
Electron density around MN A 610:

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



Electron density around MN B 601:

$2mF_o-DF_c$ (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.