



Full wwPDB X-ray Structure Validation Report ⓘ

Oct 3, 2021 – 07:54 AM EDT

PDB ID : 3I0L
Title : Crystal structure of GTB C80S/C196S/C209S + DA + UDP-Gal
Authors : Schuman, B.; Persson, M.; Landry, R.C.; Polakowski, R.; Weadge, J.T.; Seto, N.O.L.; Borisova, S.; Palcic, M.M.; Evans, S.V.
Deposited on : 2009-06-25
Resolution : 1.60 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

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 : 2.23.2
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) : 2.23.2

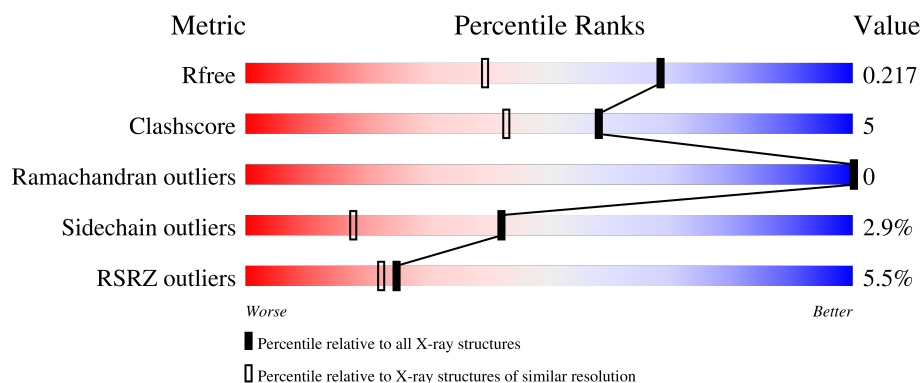
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.60 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	3398 (1.60-1.60)
Clashscore	141614	3665 (1.60-1.60)
Ramachandran outliers	138981	3564 (1.60-1.60)
Sidechain outliers	138945	3563 (1.60-1.60)
RSRZ outliers	127900	3321 (1.60-1.60)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments 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	287	<div> <div>5%</div> <div> <div></div> <div>78%</div> <div>14%</div> <div>•• 5%</div> </div> </div>
2	B	2	<div> <div>50%</div> <div>50%</div> </div>

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
4	GAL	A	359	X	-	-	-

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 2644 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 ABO glycosyltransferase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	272	Total	C	N	O	S	0	4	0
			2240	1452	381	397	10			

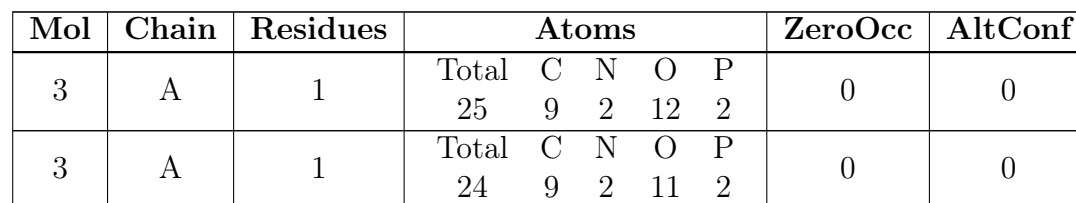
There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	80	SER	CYS	engineered mutation	UNP Q70V26
A	196	SER	CYS	engineered mutation	UNP Q70V26
A	209	SER	CYS	engineered mutation	UNP Q70V26
A	355	GLU	-	SEE REMARK 999	UNP Q70V26

- Molecule 2 is an oligosaccharide called alpha-L-fucopyranose-(1-2)-hexyl beta-D-galactopyranoside.

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace
2	B	2	Total	C	O	0	0	0
			27	18	9			

- Molecule 3 is URIDINE-5'-DIPHOSPHATE (three-letter code: UDP) (formula: C₉H₁₄N₂O₁₂P₂).



- GAL
-
- Chemical structure of Galactose (GAL) in its cyclic form (pyranose ring). The structure shows the following atoms and bonds:
- Carbons: C1(R), C2(R), C3(S), C4(R), C5(R), C6
 - Oxygen atoms: O1, O2, O3, O4, O5, O6
 - Hydroxyl groups: OH (O1, O2, O3, O4, O5)
 - Hydroxymethyl group: CH₂OH (C6, O6)
- The stereochemistry is indicated by the labels (R) and (S) for the chiral centers. The structure is shown in a chair conformation.

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			12	6	6		

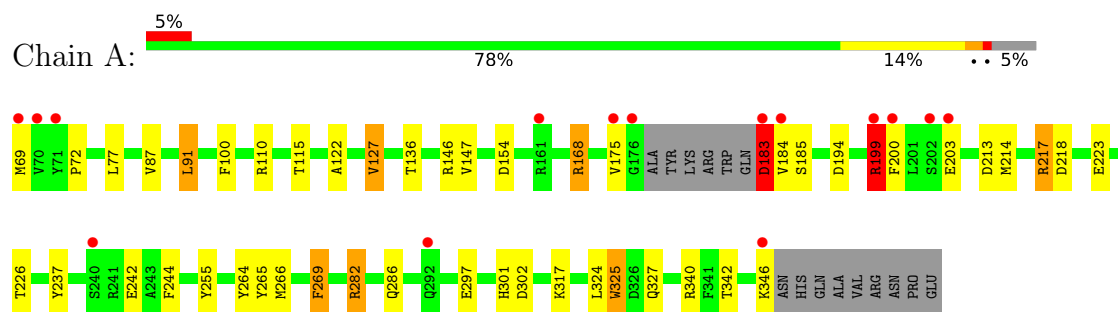
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	316	Total 316	O 316	0	0

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: ABO glycosyltransferase



- Molecule 2: alpha-L-fucopyranose-(1-2)-hexyl beta-D-galactopyranoside



4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	52.55Å 149.28Å 79.45Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	19.76 – 1.60 19.76 – 1.60	Depositor EDS
% Data completeness (in resolution range)	94.6 (19.76-1.60) 94.5 (19.76-1.60)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.93 (at 1.60Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.184 , 0.223 0.180 , 0.217	Depositor DCC
R_{free} test set	1977 reflections (5.01%)	wwPDB-VP
Wilson B-factor (Å ²)	19.1	Xtriage
Anisotropy	0.164	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.39 , 55.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	2644	wwPDB-VP
Average B, all atoms (Å ²)	19.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.98% 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

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: GAL, BHG, UDP, FUC

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z > 5$	RMSZ	$\# Z > 5$
1	A	1.41	12/2313 (0.5%)	1.31	21/3138 (0.7%)

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

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

All (12) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	237	TYR	CD2-CE2	6.84	1.49	1.39
1	A	265	TYR	CE1-CZ	6.05	1.46	1.38
1	A	194	ASP	N-CA	5.78	1.57	1.46
1	A	110	ARG	CZ-NH2	5.69	1.40	1.33
1	A	265	TYR	CE2-CZ	5.63	1.45	1.38
1	A	255	TYR	CG-CD2	5.57	1.46	1.39
1	A	286	GLN	CG-CD	5.54	1.63	1.51
1	A	325	TRP	CE3-CZ3	5.25	1.47	1.38
1	A	127	VAL	CB-CG1	-5.20	1.42	1.52
1	A	340	ARG	CZ-NH2	5.17	1.39	1.33
1	A	100	PHE	CE1-CZ	5.14	1.47	1.37
1	A	264	TYR	CG-CD2	5.04	1.45	1.39

All (21) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	282	ARG	NE-CZ-NH1	16.65	128.62	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	282	ARG	NE-CZ-NH2	-12.35	114.12	120.30
1	A	213	ASP	CB-CG-OD2	10.25	127.53	118.30
1	A	199	ARG	NE-CZ-NH2	-7.48	116.56	120.30
1	A	223	GLU	OE1-CD-OE2	-6.84	115.10	123.30
1	A	244	PHE	CB-CG-CD1	-6.84	116.02	120.80
1	A	168	ARG	NE-CZ-NH2	6.31	123.45	120.30
1	A	213	ASP	CB-CG-OD1	-6.12	112.80	118.30
1	A	154	ASP	CB-CG-OD1	5.88	123.59	118.30
1	A	136	THR	CA-CB-CG2	-5.69	104.43	112.40
1	A	244	PHE	CB-CG-CD2	5.66	124.76	120.80
1	A	302	ASP	CB-CG-OD1	5.66	123.40	118.30
1	A	168	ARG	NE-CZ-NH1	-5.66	117.47	120.30
1	A	217	ARG	NE-CZ-NH1	-5.66	117.47	120.30
1	A	91	LEU	CB-CG-CD1	-5.52	101.61	111.00
1	A	269	PHE	CB-CG-CD2	5.52	124.67	120.80
1	A	77	LEU	CB-CG-CD1	-5.36	101.89	111.00
1	A	175	VAL	CG1-CB-CG2	-5.26	102.48	110.90
1	A	218	ASP	CB-CG-OD1	5.19	122.97	118.30
1	A	214	MET	CG-SD-CE	-5.14	91.98	100.20
1	A	340	ARG	NE-CZ-NH1	5.04	122.82	120.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	183	ASP	Peptide

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	2240	0	2223	21	0
2	B	27	0	31	0	0
3	A	49	0	22	2	0
4	A	12	0	11	0	0
5	A	316	0	0	4	0
All	All	2644	0	2287	22	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 5.

All (22) 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:242:GLU:OE2	5:A:538:HOH:O	1.72	1.08
1:A:282:ARG:HD3	5:A:539:HOH:O	1.77	0.84
1:A:184:VAL:HA	5:A:522:HOH:O	1.84	0.76
1:A:217:ARG:HH22	1:A:327:GLN:HE21	1.37	0.73
1:A:297:GLU:OE1	1:A:301:HIS:HD2	1.76	0.68
1:A:217:ARG:HH12	1:A:327:GLN:HE22	1.43	0.66
1:A:217:ARG:HH12	1:A:327:GLN:NE2	1.99	0.61
3:A:358:UDP:O1A	5:A:8:HOH:O	2.17	0.57
1:A:183:ASP:CG	1:A:184:VAL:H	2.08	0.57
1:A:217:ARG:HH22	1:A:327:GLN:NE2	2.07	0.52
1:A:199:ARG:NE	1:A:199:ARG:H	2.12	0.48
1:A:325:TRP:HB3	1:A:342:THR:HG22	1.96	0.47
1:A:72:PRO:HG2	3:A:357:UDP:H5'2	1.99	0.45
1:A:183:ASP:OD1	1:A:184:VAL:N	2.49	0.45
1:A:199:ARG:HD2	1:A:200:PHE:N	2.33	0.44
1:A:266:MET:HB3	1:A:324:LEU:CD2	2.47	0.43
1:A:147:VAL:O	1:A:168:ARG:HA	2.19	0.43
1:A:226:THR:HG21	1:A:317:LYS:HB2	2.01	0.43
1:A:87[A]:VAL:CG1	1:A:91:LEU:HA	2.50	0.42
1:A:183:ASP:CG	1:A:184:VAL:N	2.72	0.41
1:A:115:THR:OG1	1:A:146:ARG:HD3	2.21	0.40
1:A:122:ALA:HB1	1:A:127:VAL:HA	2.02	0.40

There are no symmetry-related clashes.

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	272/287 (95%)	266 (98%)	6 (2%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains ⓘ

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	246/255 (96%)	238 (97%)	8 (3%)	38	14

All (8) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	69	MET
1	A	183	ASP
1	A	185[A]	SER
1	A	185[B]	SER
1	A	199	ARG
1	A	203	GLU
1	A	269	PHE
1	A	346	LYS

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

Mol	Chain	Res	Type
1	A	73	GLN
1	A	155	GLN
1	A	285	HIS
1	A	301	HIS
1	A	327	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

2 monosaccharides are modelled in this entry.

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	BHG	B	1	2	16,17,18	1.13	1 (6%)	18,21,23	1.41	3 (16%)
2	FUC	B	2	2	10,10,11	1.11	0	14,14,16	0.56	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	BHG	B	1	2	-	1/9/25/29	0/1/1/1
2	FUC	B	2	2	-	-	0/1/1/1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	1	BHG	C4-C5	2.84	1.57	1.52

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	1	BHG	C6-C5-C4	3.25	118.08	113.54
2	B	1	BHG	C1'-O1-C1	-3.12	108.67	113.84
2	B	1	BHG	C3-C4-C5	-2.27	107.83	110.77

There are no chirality outliers.

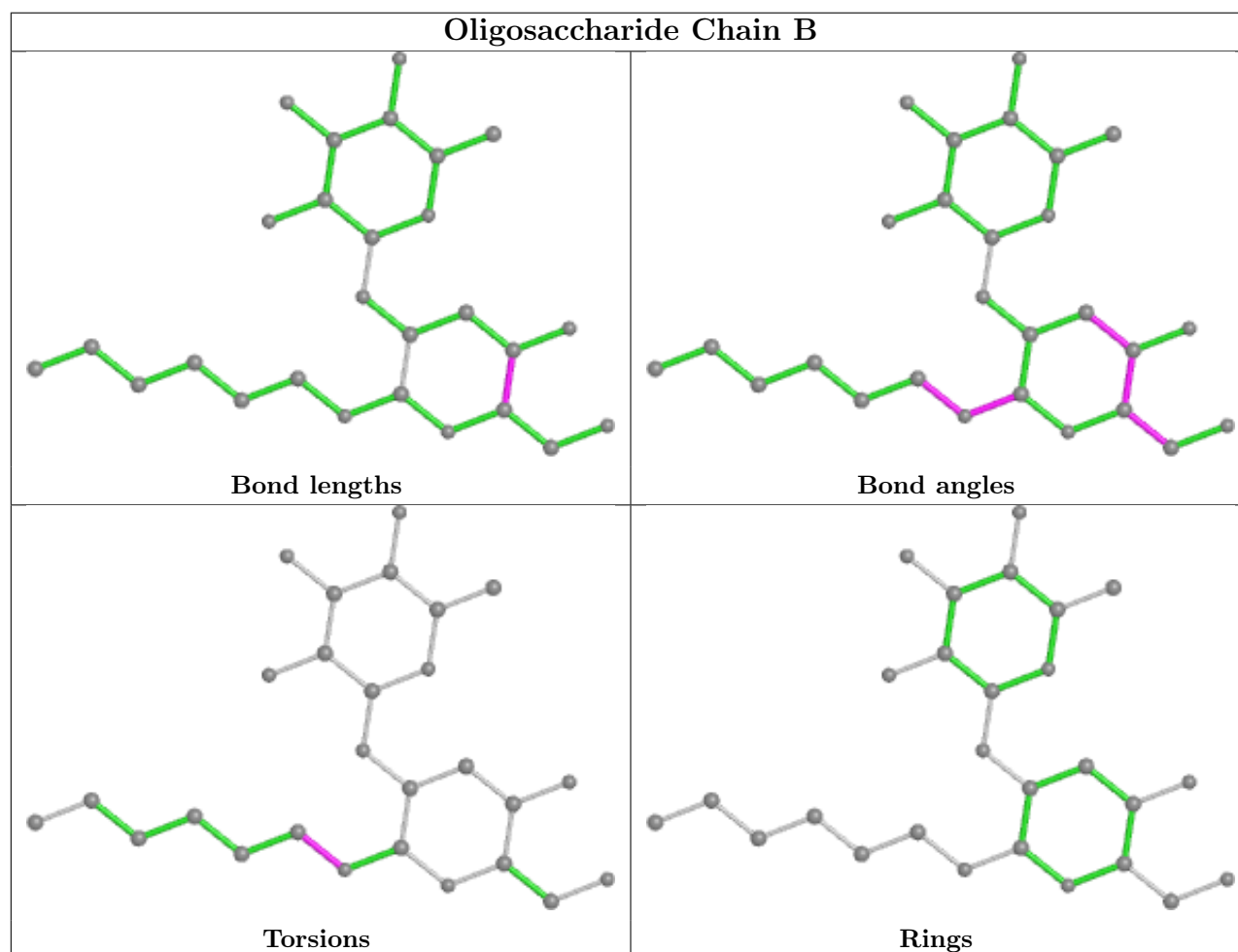
All (1) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	B	1	BHG	C2'-C1'-O1-C1

There are no ring outliers.

No monomer is involved in short contacts.

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for oligosaccharide.



5.6 Ligand geometry [i](#)

3 ligands are modelled in this entry.

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	$\# Z > 2$
3	UDP	A	358	4	18,25,26	1.91	4 (22%)	19,37,40	0.72	0
4	GAL	A	359	3	12,12,12	1.46	1 (8%)	17,17,17	0.83	1 (5%)
3	UDP	A	357	-	20,26,26	2.24	6 (30%)	25,40,40	1.10	2 (8%)

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
3	UDP	A	358	4	-	2/11/31/32	0/2/2/2
4	GAL	A	359	3	1/1/5/5	0/2/22/22	0/1/1/1
3	UDP	A	357	-	-	3/14/32/32	0/2/2/2

All (11) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	357	UDP	C6-N1	5.31	1.42	1.35
3	A	358	UDP	C4-N3	4.88	1.41	1.33
3	A	358	UDP	C6-N1	4.77	1.41	1.35
3	A	357	UDP	PB-O1B	4.74	1.65	1.50
3	A	357	UDP	C4-N3	4.31	1.40	1.33
4	A	359	GAL	O1-C1	4.13	1.52	1.39
3	A	357	UDP	PA-O1A	3.04	1.61	1.50
3	A	358	UDP	O4'-C1'	2.56	1.44	1.41
3	A	357	UDP	O4'-C1'	2.33	1.44	1.41
3	A	358	UDP	PA-O1A	2.30	1.59	1.50
3	A	357	UDP	PB-O2B	2.18	1.63	1.54

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	357	UDP	O3B-PB-O3A	3.30	115.70	104.64
3	A	357	UDP	PA-O3A-PB	-2.38	124.66	132.83
4	A	359	GAL	O1-C1-C2	2.07	114.87	109.03

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
4	A	359	GAL	C1

All (5) torsion outliers are listed below:

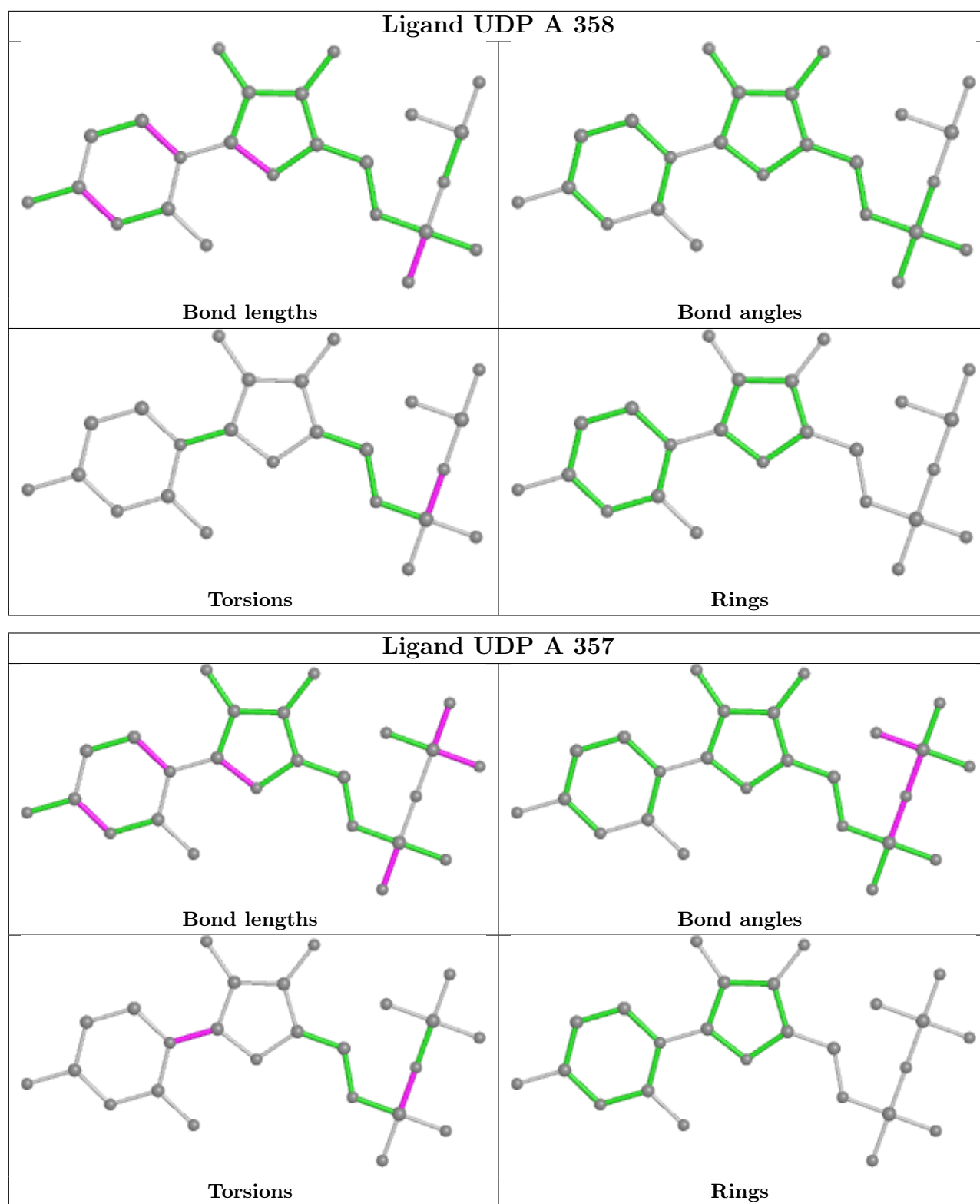
Mol	Chain	Res	Type	Atoms
3	A	357	UDP	C2'-C1'-N1-C6
3	A	357	UDP	O4'-C1'-N1-C6
3	A	358	UDP	PB-O3A-PA-O1A
3	A	357	UDP	PB-O3A-PA-O2A
3	A	358	UDP	PB-O3A-PA-O2A

There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	358	UDP	1	0
3	A	357	UDP	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 ⓘ

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, 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	272/287 (94%)	0.21	15 (5%)	25 22	9, 15, 34, 51	0

All (15) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	346	LYS	6.0
1	A	69	MET	5.4
1	A	176	GLY	4.0
1	A	70	VAL	3.7
1	A	203	GLU	3.7
1	A	240	SER	3.2
1	A	199	ARG	3.1
1	A	202	SER	3.0
1	A	161	ARG	3.0
1	A	71	TYR	2.9
1	A	184	VAL	2.9
1	A	200	PHE	2.7
1	A	183	ASP	2.2
1	A	175	VAL	2.1
1	A	292	GLN	2.1

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

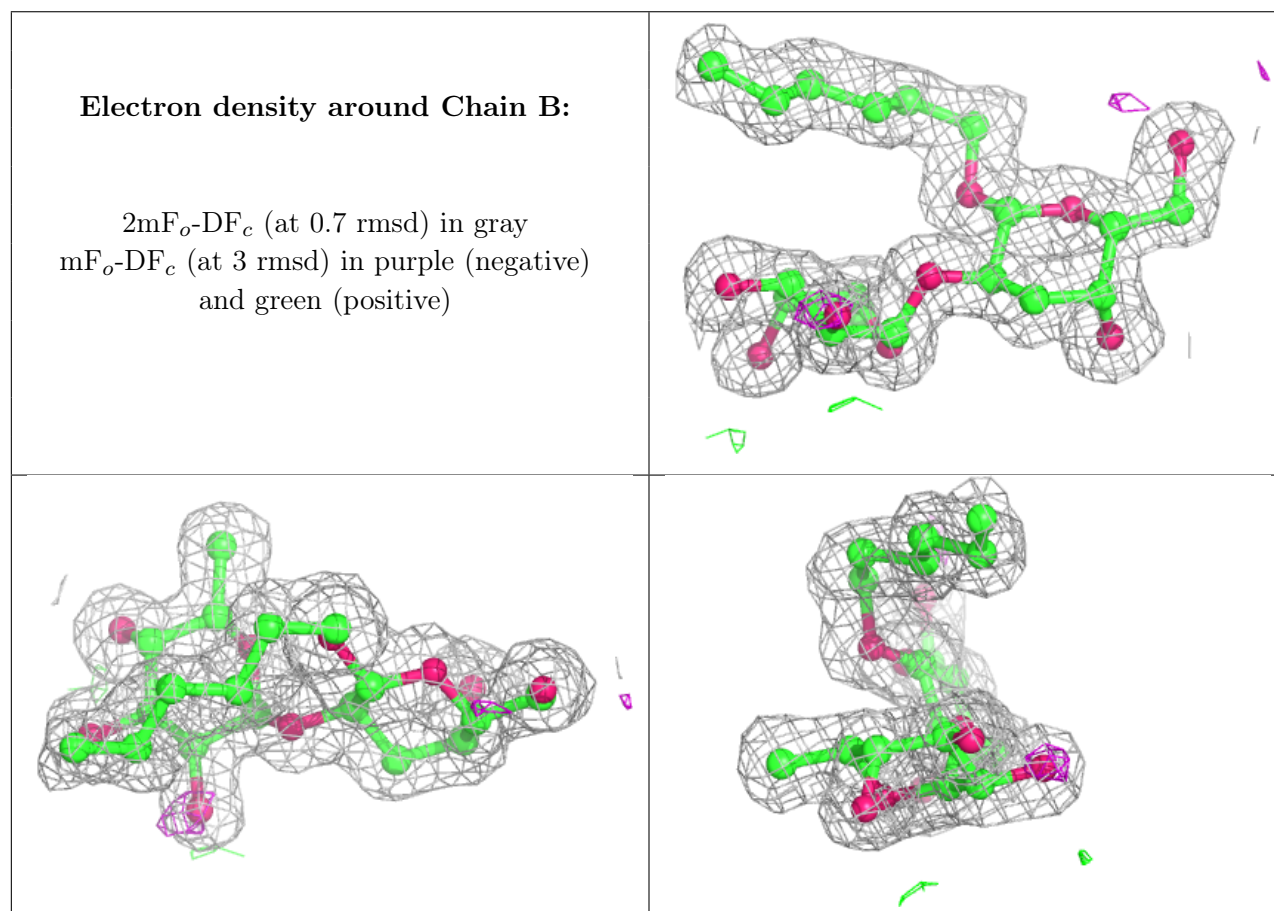
There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates [i](#)

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
2	FUC	B	2	10/11	0.88	0.10	14,15,16,17	0
2	BHG	B	1	17/18	0.94	0.08	12,15,23,27	0

The following is a graphical depiction of the model fit to experimental electron density for oligosaccharide. Each fit is shown from different orientation to approximate a three-dimensional view.



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.

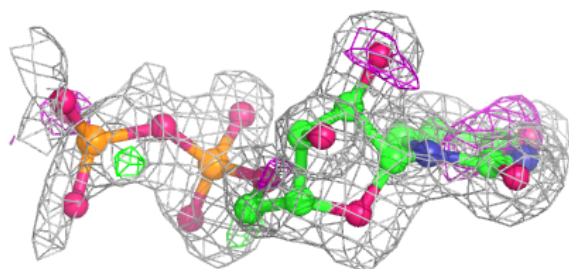
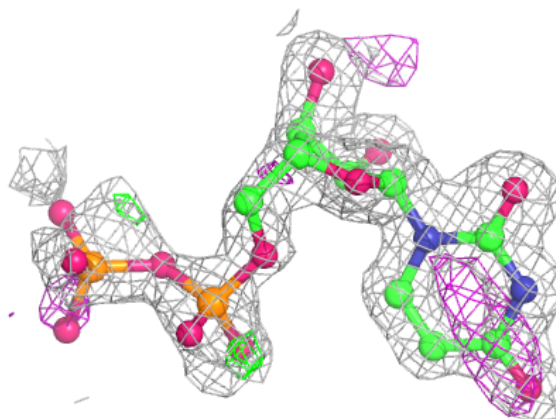
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
3	UDP	A	357	25/25	0.81	0.21	19,21,37,39	8
4	GAL	A	359	12/12	0.88	0.15	19,21,25,26	0
3	UDP	A	358	24/25	0.92	0.14	14,20,22,25	0

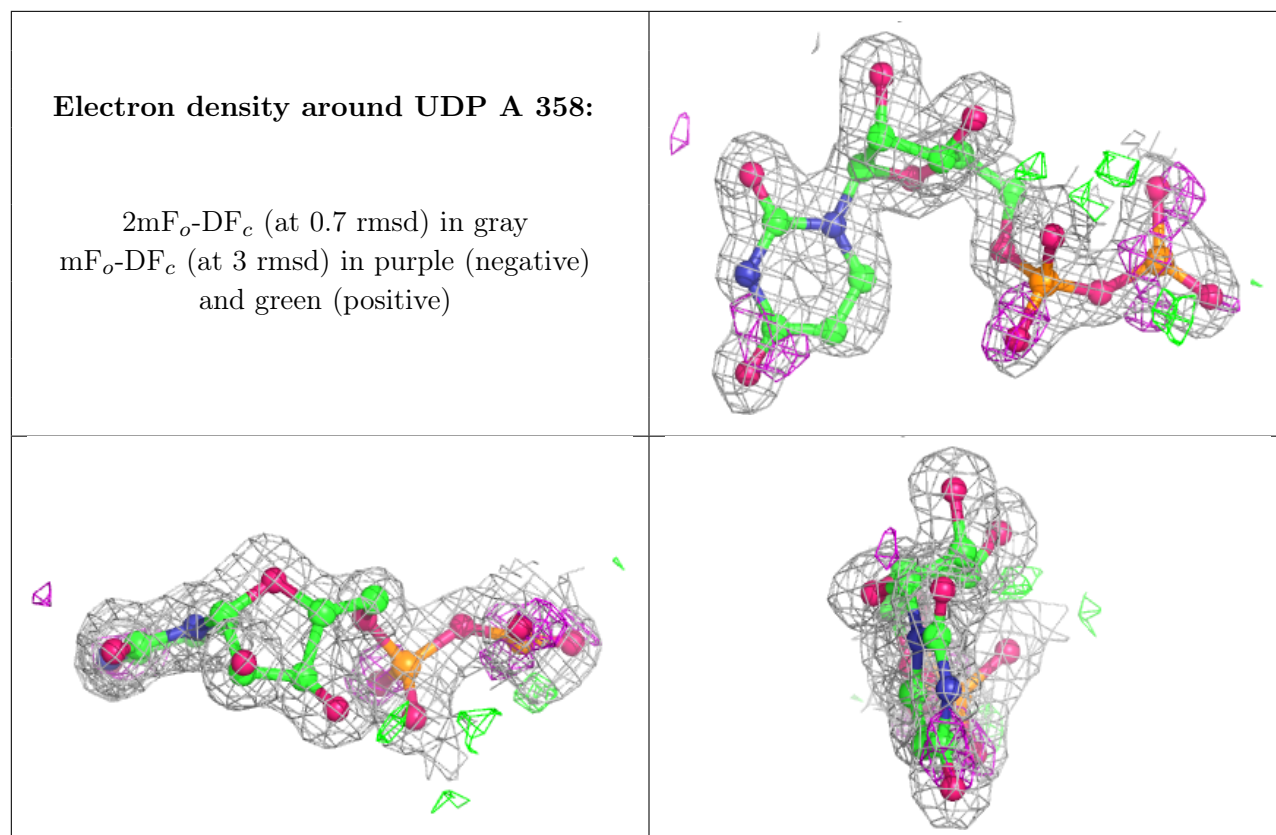
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers

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 UDP A 357:

$2mF_o-DF_c$ (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.