



# Full wwPDB X-ray Structure Validation Report ⓘ

Aug 10, 2020 – 09:53 AM BST

PDB ID : 6NEA  
Title : Human Acetylcholinesterase in complex with reactivator, HLo7  
Authors : Bester, S.M.; McGuire, J.; Height, J.J.; Pegan, S.D.  
Deposited on : 2018-12-17  
Resolution : 2.42 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.

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

MolProbity : 4.02b-467  
Mogul : 1.8.5 (274361), CSD as541be (2020)  
Xtriage (Phenix) : 1.13  
EDS : 2.13.1  
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

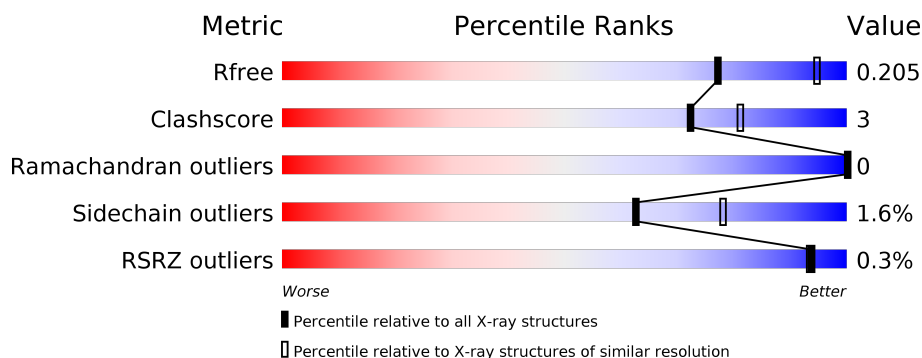
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.42 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	4647 (2.44-2.40)
Clashscore	141614	5161 (2.44-2.40)
Ramachandran outliers	138981	5073 (2.44-2.40)
Sidechain outliers	138945	5074 (2.44-2.40)
RSRZ outliers	127900	4543 (2.44-2.40)

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  $\geq 3$ , 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	542	<div> <div>91%</div> <div>8%</div> <div>.</div> </div>
1	B	542	<div> <div>88%</div> <div>9%</div> <div>.</div> </div>
2	C	3	<div> <div>67%</div> <div>33%</div> </div>
2	D	3	<div> <div>67%</div> <div>33%</div> </div>

## 2 Entry composition [i](#)

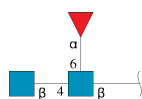
There are 6 unique types of molecules in this entry. The entry contains 9103 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Acetylcholinesterase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	533	Total	C	N	O	S	0	1	0
			4150	2664	727	746	13			
1	B	528	Total	C	N	O	S	0	3	0
			4130	2654	724	739	13			

- Molecule 2 is an oligosaccharide called 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-[alpha-L-fucopyranose-(1-6)]2-acetamido-2-deoxy-beta-D-glucopyranose.



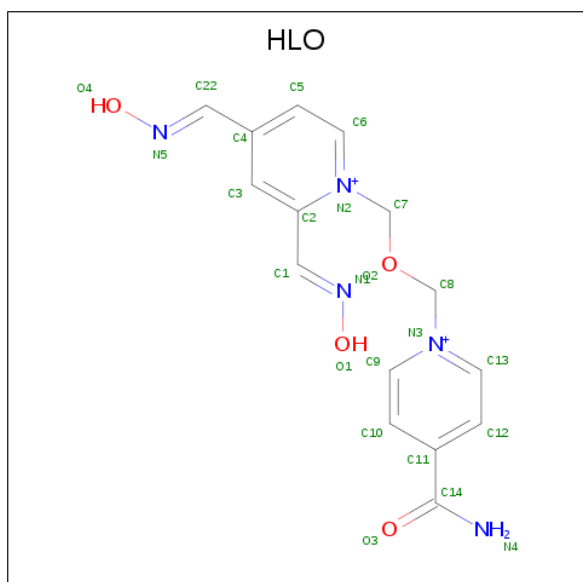
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	C	3	Total	C	N	O	0	0	0
			38	22	2	14			
2	D	3	Total	C	N	O	0	0	0
			38	22	2	14			

- Molecule 3 is 2-acetamido-2-deoxy-beta-D-glucopyranose (three-letter code: NAG) (formula: C<sub>8</sub>H<sub>15</sub>NO<sub>6</sub>).



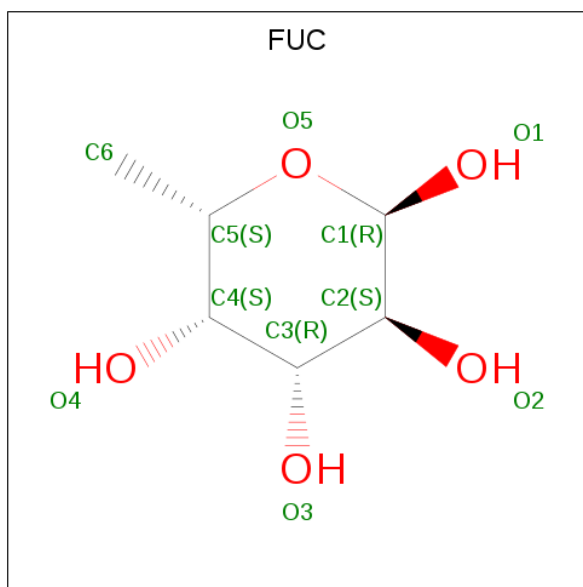
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			14	8	1	5		
3	A	1	Total	C	N	O	0	0
			14	8	1	5		
3	A	1	Total	C	N	O	0	0
			14	8	1	5		
3	B	1	Total	C	N	O	0	0
			14	8	1	5		

- Molecule 4 is 1-[(2,4-BIS[(E)-(HYDROXYIMINO)METHYL]PYRIDINIUM-1-YL)METHOXY]METHYL-4-CARBAMOYLPYRIDINIUM (three-letter code: HLO) (formula:  $C_{15}H_{17}N_5O_4$ ).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	1	Total	C	N	O	0	0
			24	15	5	4		
4	B	1	Total	C	N	O	0	0
			24	15	5	4		

- Molecule 5 is alpha-L-fucopyranose (three-letter code: FUC) (formula: C<sub>6</sub>H<sub>12</sub>O<sub>5</sub>).



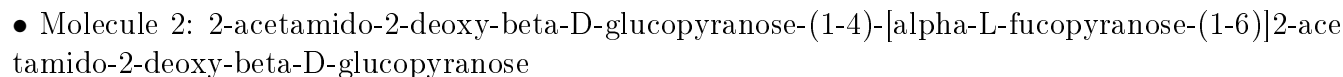
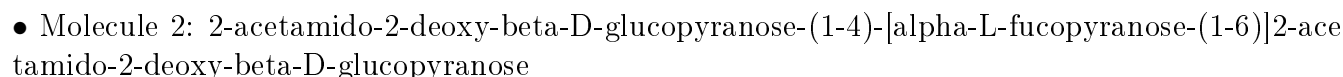
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	C	O	0	0
			10	6	4		

- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	286	Total	O	0	0
			286	286		
6	B	347	Total	O	0	0
			347	347		



- Molecule 1: Acetylcholinesterase



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 31 2 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	104.73Å 104.73Å 324.25Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	41.25 – 2.42 46.43 – 2.42	Depositor EDS
% Data completeness (in resolution range)	99.6 (41.25-2.42) 99.6 (46.43-2.42)	Depositor EDS
$R_{merge}$	0.11	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	3.31 (at 2.42Å)	Xtriage
Refinement program	PHENIX (dev_3311: ???)	Depositor
R, $R_{free}$	0.176 , 0.205 0.176 , 0.205	Depositor DCC
$R_{free}$ test set	3964 reflections (4.98%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	43.9	Xtriage
Anisotropy	0.531	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.32 , 43.3	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	0.016 for -h,-k,l	Xtriage
$F_o, F_c$ correlation	0.96	EDS
Total number of atoms	9103	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	49.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.60% 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

### 5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: HLO, NAG, 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	0.27	0/4279	0.46	0/5847
1	B	0.26	0/4265	0.44	0/5828
All	All	0.27	0/8544	0.45	0/11675

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

### 5.2 Too-close contacts

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	4150	0	4033	25	0
1	B	4130	0	4021	32	0
2	C	38	0	34	1	0
2	D	38	0	34	0	0
3	A	42	0	39	1	0
3	B	14	0	13	0	0
4	A	24	0	17	3	0
4	B	24	0	17	2	0
5	A	10	0	10	0	0
6	A	286	0	0	1	0
6	B	347	0	0	2	0
All	All	9103	0	8218	57	0



The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 3.

All (57) 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:117:TRP:HB2	1:B:200:PHE:CE2	2.20	0.77
1:B:203:SER:OG	6:B:701:HOH:O	2.13	0.66
1:A:113:PRO:HG3	1:A:485:ARG:HG2	1.78	0.64
4:B:605:HLO:O1	4:B:605:HLO:H7C1	2.03	0.59
1:B:285:GLU:HG2	4:B:605:HLO:N4	2.19	0.57
1:B:117:TRP:CE3	1:B:200:PHE:CE1	2.93	0.57
1:A:285:GLU:OE2	4:A:605:HLO:N4	2.36	0.56
1:B:117:TRP:CE3	1:B:200:PHE:CZ	2.95	0.54
1:A:203:SER:HB2	1:A:447:HIS:NE2	2.22	0.53
1:A:309:SER:OG	6:A:701:HOH:O	2.18	0.53
1:A:68:VAL:HG23	1:A:90:ARG:HB2	1.89	0.53
1:B:313:GLU:OE1	6:B:702:HOH:O	2.17	0.53
1:A:203:SER:OG	1:A:204:ALA:N	2.46	0.49
1:B:113:PRO:HG3	1:B:485:ARG:HG2	1.94	0.49
1:A:25:PRO:O	1:B:325:GLN:NE2	2.46	0.49
1:B:68:VAL:HG23	1:B:90:ARG:HB2	1.95	0.49
1:A:463:ARG:HA	3:A:604:NAG:H82	1.93	0.48
1:B:117:TRP:HB2	1:B:200:PHE:HE2	1.72	0.48
1:A:347:SER:HB2	2:C:1:NAG:H62	1.96	0.48
1:B:103:THR:HG21	1:B:190:PHE:HB3	1.95	0.48
1:B:227:LEU:HB2	1:B:328:VAL:HG12	1.95	0.48
1:A:258:PRO:HD3	1:A:269:LEU:HD23	1.96	0.47
1:A:200:PHE:CB	1:A:226:VAL:HB	2.44	0.47
1:B:203:SER:HB2	1:B:447:HIS:NE2	2.30	0.47
1:A:512:SER:OG	1:A:519:GLU:HB3	2.16	0.46
1:A:265:ASN:HD22	1:A:268:GLU:CD	2.19	0.45
1:A:458:PRO:HA	1:A:465:TYR:CD1	2.51	0.45
1:A:224:ARG:HG3	1:A:224:ARG:HH11	1.82	0.45
1:B:329:GLY:HA3	1:B:428:TYR:CE2	2.52	0.45
1:A:224:ARG:HG3	1:A:224:ARG:NH1	2.32	0.44
1:B:286:TRP:CH2	1:B:297:PHE:HA	2.53	0.43
1:A:32:PHE:HB3	1:A:35:ILE:HD11	2.00	0.43
1:A:205:GLY:O	1:A:209:VAL:HG23	2.18	0.43
1:B:404:ASP:HA	1:B:408:VAL:HB	2.00	0.43
1:A:161:LEU:HD11	1:A:269:LEU:HD12	2.00	0.43
1:B:329:GLY:HA3	1:B:428:TYR:CZ	2.53	0.43
1:A:25:PRO:HG3	1:A:461:PRO:HD3	2.00	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:285:GLU:HG3	1:B:286:TRP:HE3	1.84	0.42
1:B:433:ARG:NH2	1:B:439:TRP:O	2.50	0.42
1:B:328:VAL:O	1:B:427:ALA:HA	2.21	0.41
1:B:45:ARG:NE	1:B:51:GLU:OE2	2.52	0.41
1:B:453:PHE:HB3	1:B:476:LEU:HD12	2.02	0.41
1:B:25:PRO:HG3	1:B:461:PRO:HD3	2.03	0.41
1:B:460:ASP:HB3	1:B:463:ARG:HG3	2.02	0.41
1:B:413:GLN:O	1:B:417:ARG:HG2	2.21	0.41
1:A:166:GLU:OE2	1:A:274:ARG:NH2	2.48	0.41
1:B:13:ARG:CZ	1:B:185:GLU:HG2	2.50	0.41
1:B:305:GLY:HA2	1:B:309:SER:HA	2.03	0.41
1:B:200:PHE:HZ	1:B:455:PHE:CZ	2.39	0.41
1:B:158:PHE:HE2	1:B:173:LEU:HG	1.86	0.41
1:A:102:TRP:HB3	1:A:139:VAL:HG21	2.03	0.41
1:A:468:GLU:H	1:A:468:GLU:CD	2.24	0.41
1:B:458:PRO:HA	1:B:465:TYR:CD1	2.56	0.41
1:A:124:TYR:CE1	4:A:605:HLO:H9	2.56	0.40
1:B:226:VAL:HG22	1:B:327:LEU:HB3	2.02	0.40
1:A:74:ASP:HB2	4:A:605:HLO:O1	2.21	0.40
1:B:293:SER:HA	1:B:365:VAL:HG13	2.04	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	528/542 (97%)	511 (97%)	17 (3%)	0	100	100
1	B	525/542 (97%)	509 (97%)	16 (3%)	0	100	100
All	All	1053/1084 (97%)	1020 (97%)	33 (3%)	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	432/437 (99%)	425 (98%)	7 (2%)	62	78
1	B	431/437 (99%)	424 (98%)	7 (2%)	62	78
All	All	863/874 (99%)	849 (98%)	14 (2%)	62	78

All (14) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	70	TYR
1	A	146	LEU
1	A	165	ARG
1	A	285	GLU
1	A	286	TRP
1	A	295	PHE
1	A	337	TYR
1	B	70	TYR
1	B	165	ARG
1	B	200	PHE
1	B	286	TRP
1	B	295	PHE
1	B	337	TYR
1	B	462	SER

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	265	ASN

### 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 ⓘ

6 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	NAG	C	1	1,2	14,14,15	0.26	0	17,19,21	0.78	1 (5%)
2	NAG	C	2	2	14,14,15	0.39	0	17,19,21	0.35	0
2	FUC	C	3	2	10,10,11	0.66	0	14,14,16	0.79	0
2	NAG	D	1	1,2	14,14,15	0.36	0	17,19,21	0.51	0
2	NAG	D	2	2	14,14,15	0.57	0	17,19,21	0.57	0
2	FUC	D	3	2	10,10,11	0.96	0	14,14,16	1.12	2 (14%)

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	NAG	C	1	1,2	-	2/6/23/26	0/1/1/1
2	NAG	C	2	2	-	0/6/23/26	0/1/1/1
2	FUC	C	3	2	-	-	0/1/1/1
2	NAG	D	1	1,2	-	4/6/23/26	0/1/1/1
2	NAG	D	2	2	-	0/6/23/26	0/1/1/1
2	FUC	D	3	2	-	-	0/1/1/1

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	3	FUC	O5-C5-C4	2.85	114.64	109.52

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	1	NAG	C1-O5-C5	2.68	115.82	112.19
2	D	3	FUC	C1-O5-C5	2.19	117.75	112.78

There are no chirality outliers.

All (6) torsion outliers are listed below:

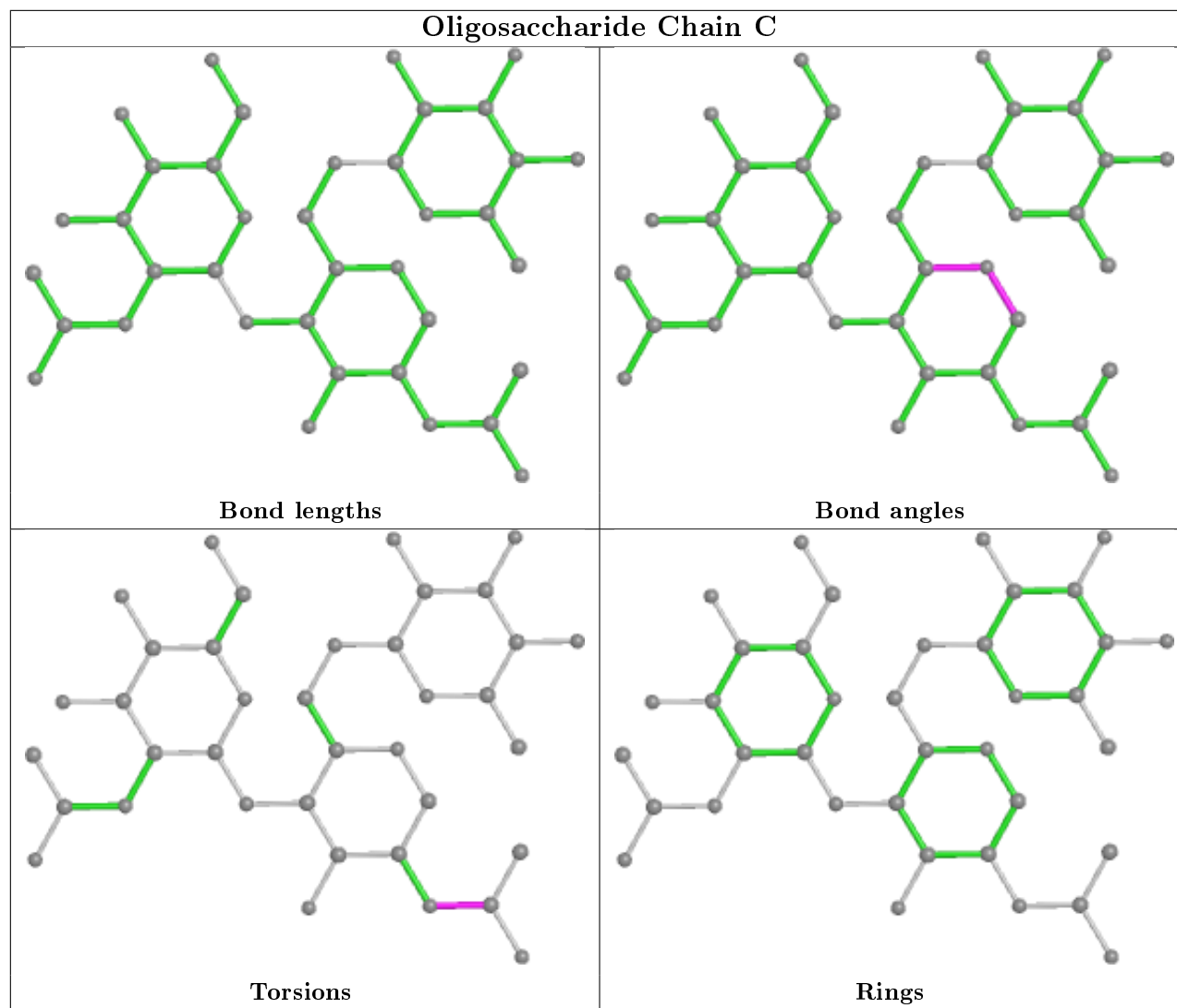
Mol	Chain	Res	Type	Atoms
2	D	1	NAG	O5-C5-C6-O6
2	D	1	NAG	C4-C5-C6-O6
2	C	1	NAG	C8-C7-N2-C2
2	C	1	NAG	O7-C7-N2-C2
2	D	1	NAG	C8-C7-N2-C2
2	D	1	NAG	O7-C7-N2-C2

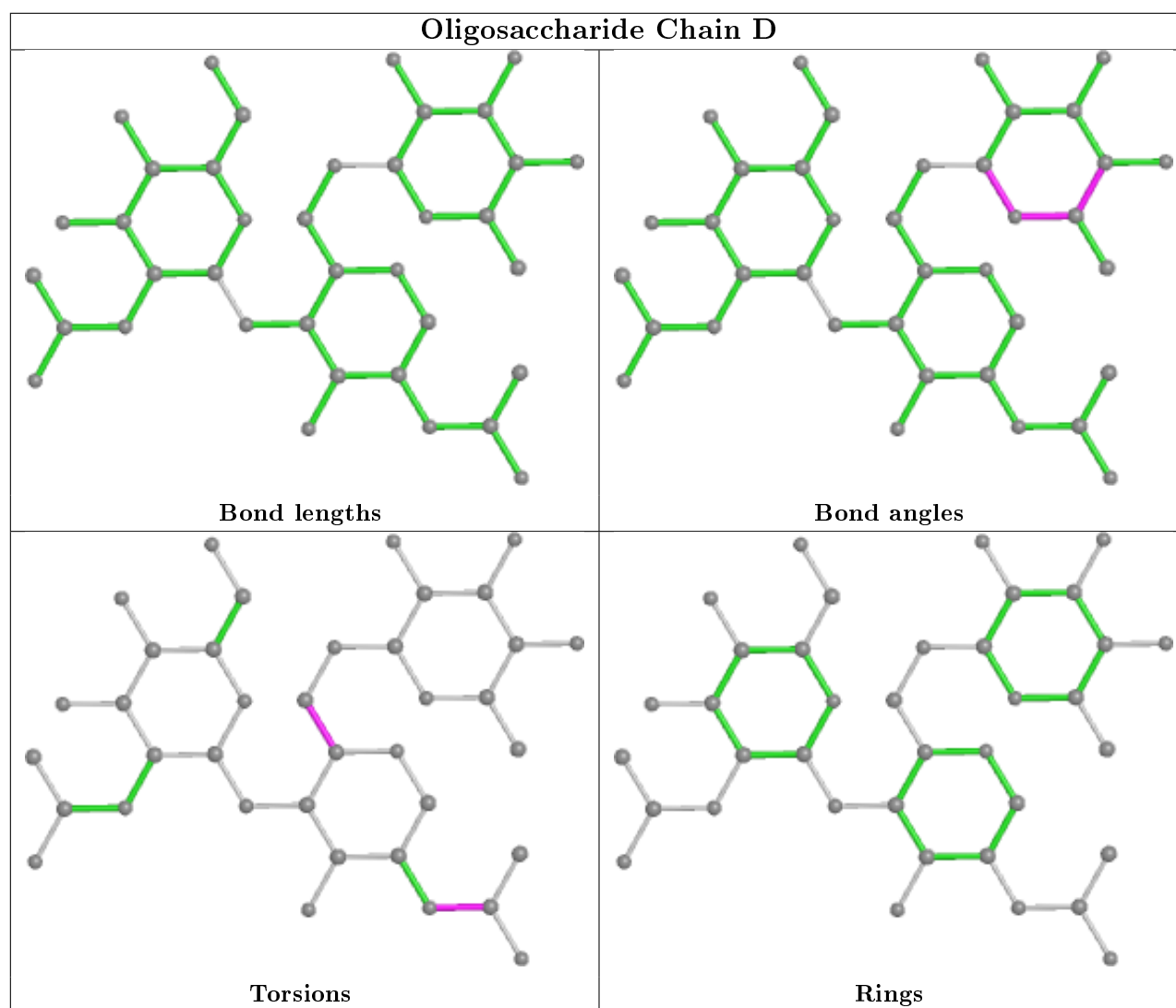
There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	1	NAG	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 oligosaccharide.





## 5.6 Ligand geometry [i](#)

7 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	NAG	B	604	1	14,14,15	0.58	0	17,19,21	0.53	0
4	HLO	A	605	-	21,25,25	2.02	6 (28%)	22,32,32	1.72	3 (13%)
3	NAG	A	607	-	14,14,15	0.26	0	17,19,21	0.40	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
5	FUC	A	606	-	10,10,11	1.32	1 (10%)	14,14,16	1.35	2 (14%)
3	NAG	A	604	1	14,14,15	0.65	1 (7%)	17,19,21	0.64	0
4	HLO	B	605	-	21,25,25	2.07	7 (33%)	22,32,32	3.10	4 (18%)
3	NAG	A	608	-	14,14,15	0.82	1 (7%)	17,19,21	0.59	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
3	NAG	B	604	1	-	2/6/23/26	0/1/1/1
4	HLO	A	605	-	-	4/15/16/16	0/2/2/2
3	NAG	A	607	-	-	2/6/23/26	0/1/1/1
5	FUC	A	606	-	-	-	0/1/1/1
3	NAG	A	604	1	-	2/6/23/26	0/1/1/1
4	HLO	B	605	-	-	4/15/16/16	0/2/2/2
3	NAG	A	608	-	-	4/6/23/26	0/1/1/1

All (16) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	B	605	HLO	C14-N4	6.73	1.45	1.33
4	A	605	HLO	C14-N4	6.42	1.45	1.33
4	B	605	HLO	C2-C1	2.91	1.54	1.46
3	A	608	NAG	O5-C1	-2.90	1.39	1.43
4	A	605	HLO	C4-C22	2.79	1.53	1.47
5	A	606	FUC	O5-C1	-2.68	1.39	1.43
4	A	605	HLO	O3-C14	-2.65	1.19	1.24
4	B	605	HLO	C4-C22	2.65	1.52	1.47
4	B	605	HLO	O3-C14	-2.60	1.19	1.24
4	B	605	HLO	C3-C4	-2.51	1.35	1.39
4	A	605	HLO	C3-C4	-2.35	1.36	1.39
4	A	605	HLO	C2-C1	2.32	1.53	1.46
3	A	604	NAG	O5-C1	-2.19	1.40	1.43
4	A	605	HLO	C12-C11	-2.10	1.35	1.39
4	B	605	HLO	C5-C4	-2.08	1.35	1.39
4	B	605	HLO	C10-C11	-2.07	1.35	1.39

All (9) bond angle outliers are listed below:



Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	605	HLO	O1-N1-C1	9.49	128.62	111.86
4	B	605	HLO	C2-C1-N1	8.76	134.69	117.75
4	A	605	HLO	O1-N1-C1	5.39	121.37	111.86
4	B	605	HLO	O4-N5-C22	4.61	120.01	111.86
4	A	605	HLO	O4-N5-C22	3.96	118.86	111.86
4	B	605	HLO	C3-C2-C1	-3.21	114.20	121.26
5	A	606	FUC	O5-C5-C4	2.65	114.27	109.52
5	A	606	FUC	C3-C4-C5	2.52	113.69	109.77
4	A	605	HLO	C5-C6-N2	-2.01	119.68	121.28

There are no chirality outliers.

All (18) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	605	HLO	O2-C8-N3-C13
4	A	605	HLO	N1-C1-C2-N2
4	A	605	HLO	N1-C1-C2-C3
3	A	607	NAG	C3-C2-N2-C7
4	B	605	HLO	O2-C7-N2-C6
4	B	605	HLO	N1-C1-C2-N2
3	A	608	NAG	O5-C5-C6-O6
3	A	604	NAG	O5-C5-C6-O6
3	A	608	NAG	C4-C5-C6-O6
3	A	608	NAG	C8-C7-N2-C2
3	A	608	NAG	O7-C7-N2-C2
3	A	604	NAG	C4-C5-C6-O6
4	B	605	HLO	N3-C8-O2-C7
4	A	605	HLO	N3-C8-O2-C7
4	B	605	HLO	C2-C1-N1-O1
3	B	604	NAG	C3-C2-N2-C7
3	B	604	NAG	C1-C2-N2-C7
3	A	607	NAG	C1-C2-N2-C7

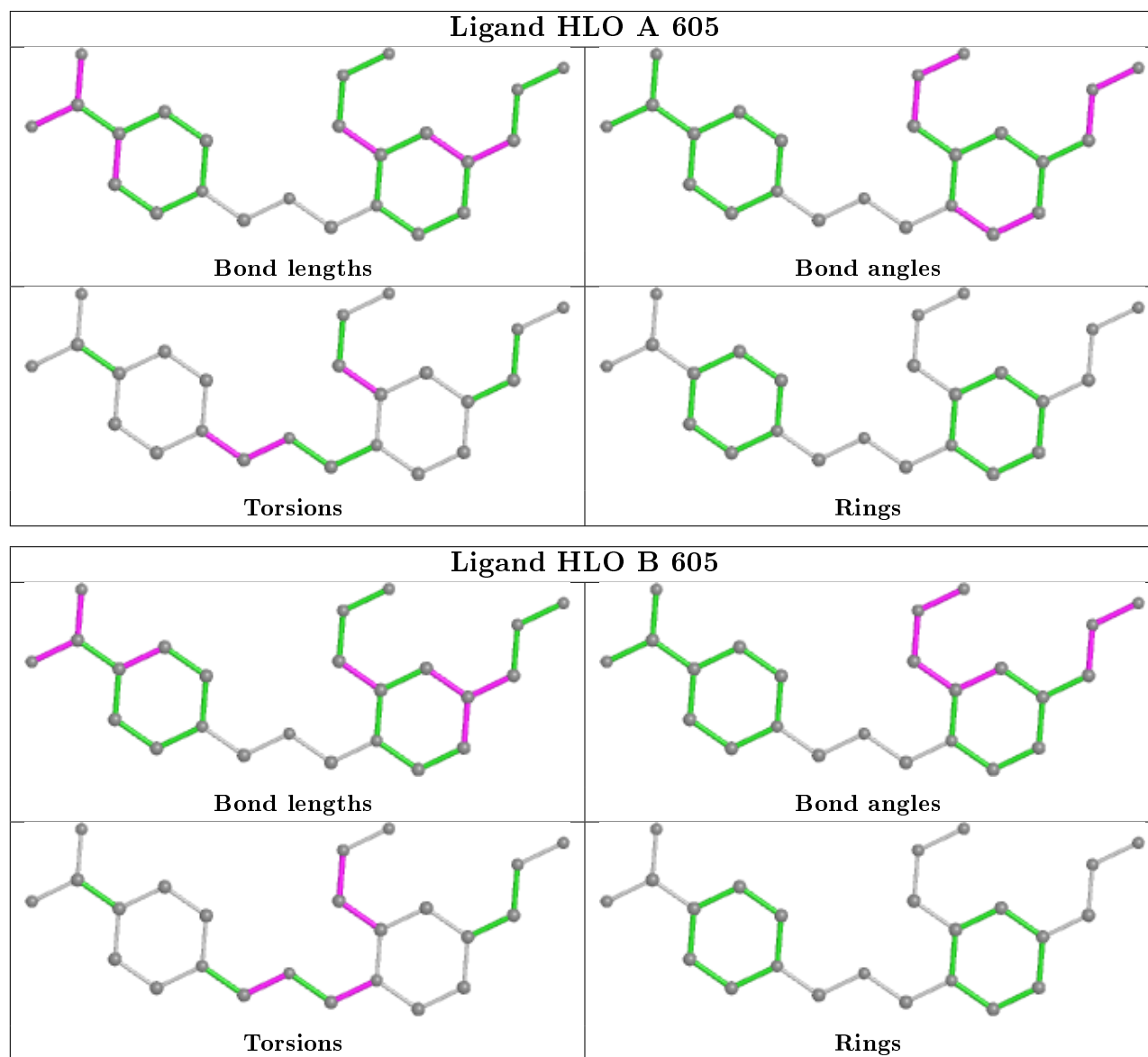
There are no ring outliers.

3 monomers are involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	605	HLO	3	0
3	A	604	NAG	1	0
4	B	605	HLO	2	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 ⓘ

There are no chain breaks in this entry.

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

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

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95<sup>th</sup> percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å <sup>2</sup> )	Q<0.9
1	A	533/542 (98%)	-0.44	2 (0%) 92 91	32, 47, 70, 113	0
1	B	528/542 (97%)	-0.52	1 (0%) 95 95	30, 44, 64, 93	0
All	All	1061/1084 (97%)	-0.48	3 (0%) 94 93	30, 45, 68, 113	0

All (3) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	264	GLY	4.3
1	A	259	PRO	3.0
1	B	542	ALA	2.1

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

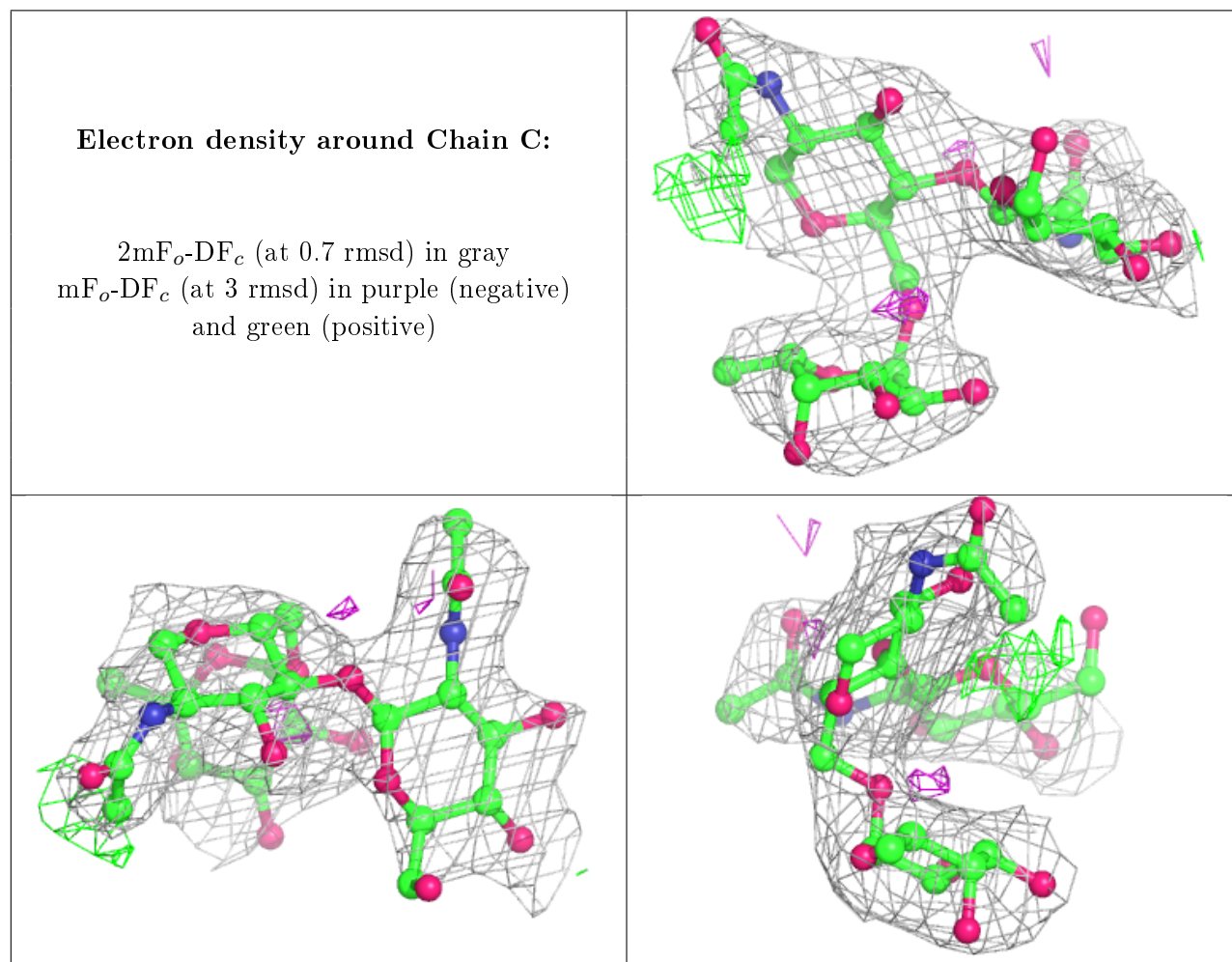
There are no non-standard protein/DNA/RNA residues in this entry.

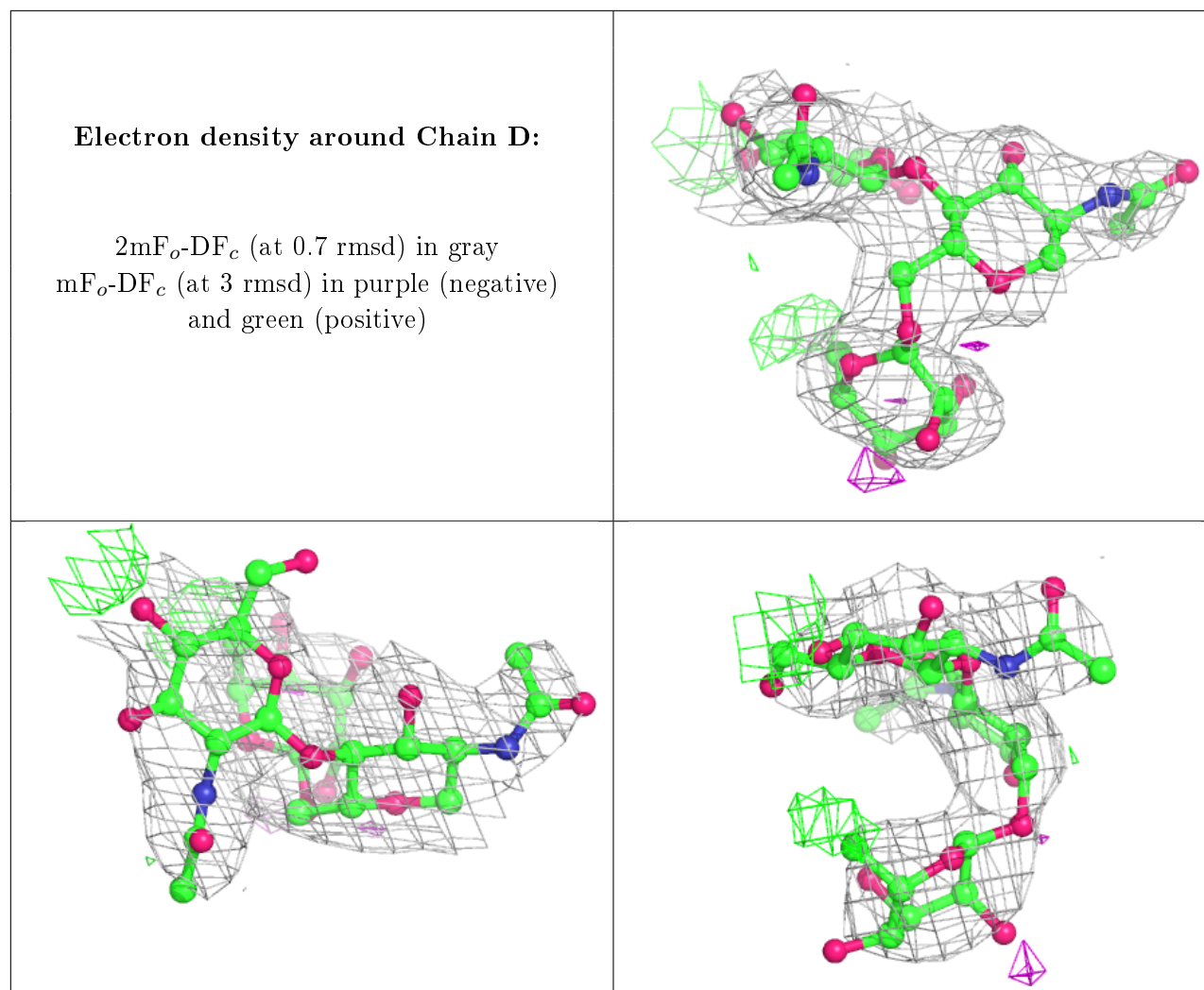
### 6.3 Carbohydrates [i](#)

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å <sup>2</sup> )	Q<0.9
2	NAG	C	2	14/15	0.85	0.34	63,103,113,118	0
2	NAG	D	2	14/15	0.87	0.34	84,113,126,128	0
2	FUC	D	3	10/11	0.89	0.40	93,105,109,111	0
2	NAG	D	1	14/15	0.92	0.29	68,90,95,100	0
2	NAG	C	1	14/15	0.92	0.27	75,84,96,102	0
2	FUC	C	3	10/11	0.94	0.32	81,97,105,107	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.





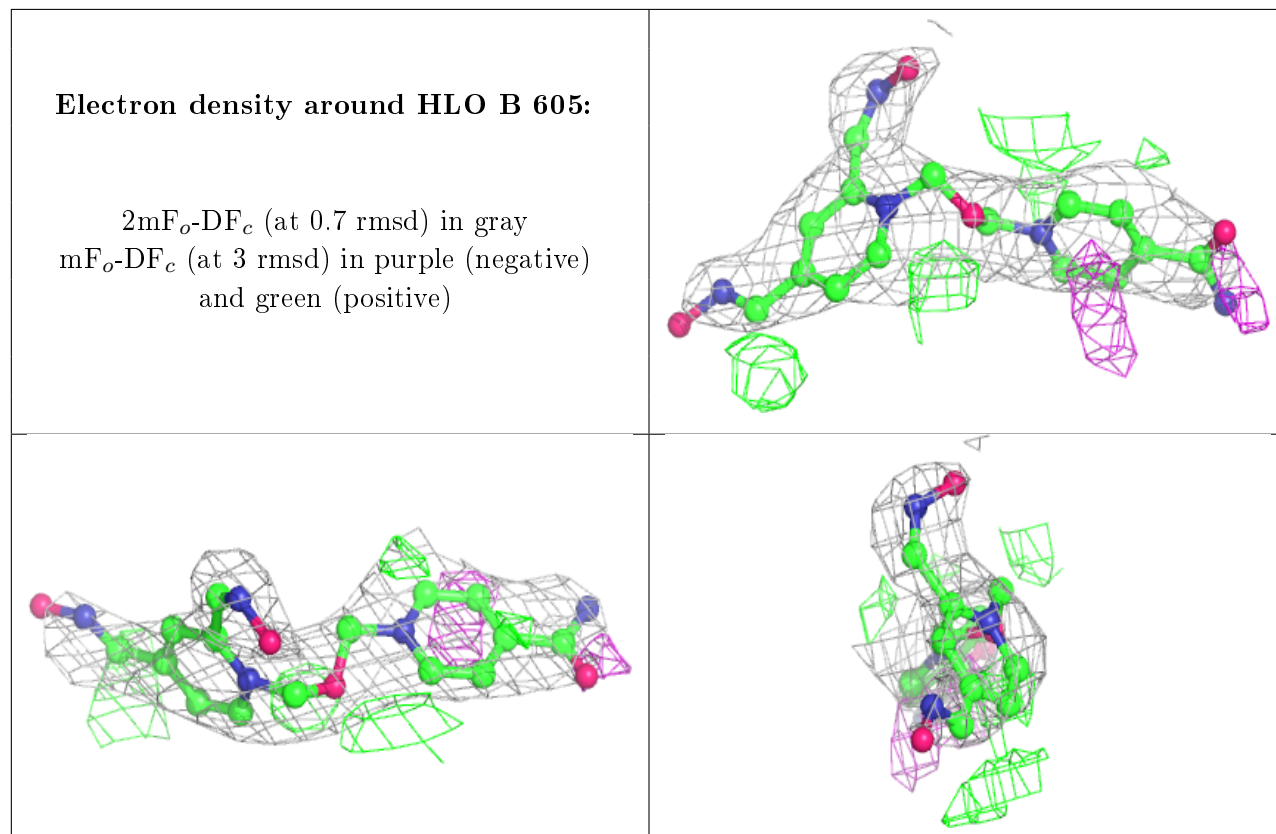
## 6.4 Ligands [i](#)

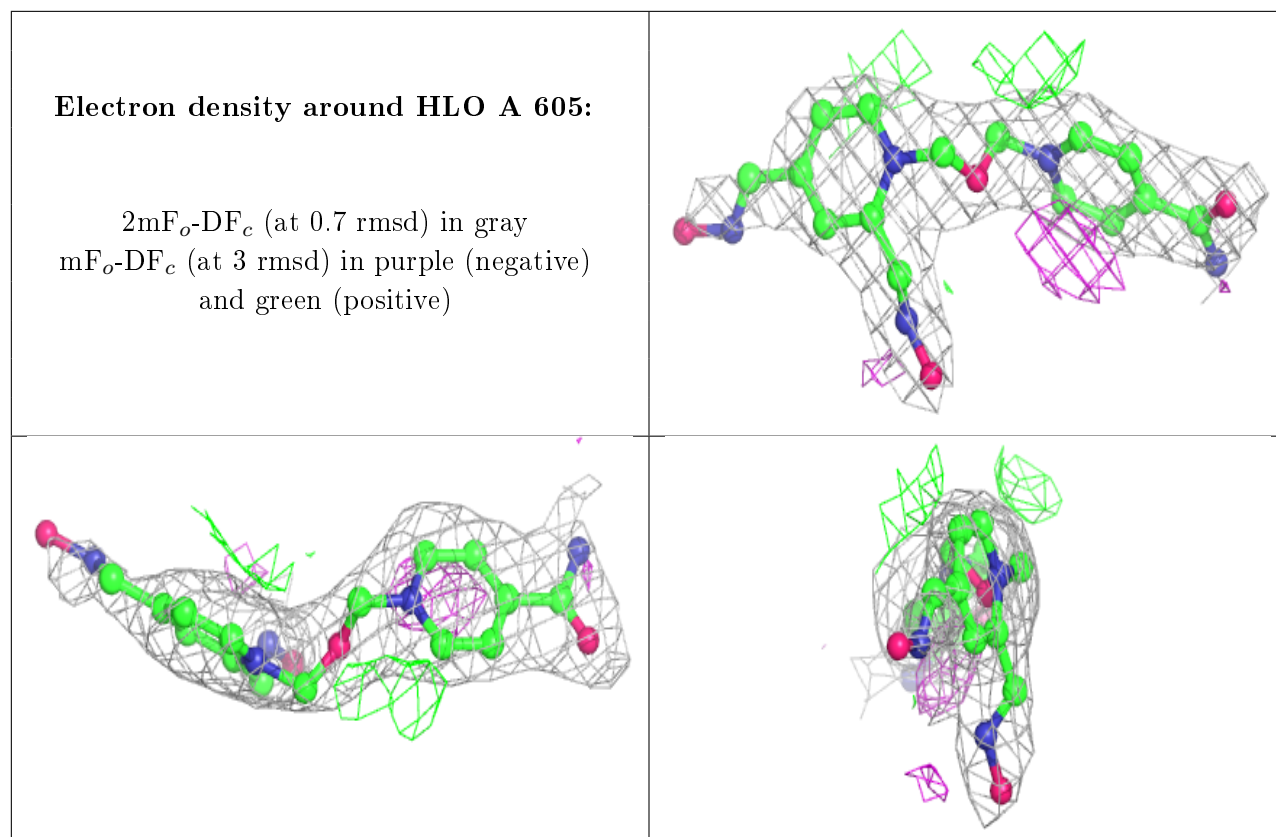
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å <sup>2</sup> )	Q<0.9
3	NAG	A	608	14/15	0.64	0.36	119,130,136,140	0
5	FUC	A	606	10/11	0.66	0.39	121,129,136,138	0
3	NAG	A	607	14/15	0.79	0.32	97,116,128,134	0
3	NAG	B	604	14/15	0.81	0.40	118,124,130,131	0
3	NAG	A	604	14/15	0.83	0.39	94,112,123,126	0
4	HLO	B	605	24/24	0.88	0.20	59,74,93,95	0
4	HLO	A	605	24/24	0.95	0.15	42,60,76,83	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.





## 6.5 Other polymers [i](#)

There are no such residues in this entry.