



wwPDB X-ray Structure Validation Summary Report ⓘ

Oct 5, 2024 – 05:27 PM EDT

PDB ID : 6MHR
Title : Structure of the human 4-1BB / Urelumab Fab complex
Authors : Kimberlin, C.R.; Chin, S.M.; Roe-Zurz, Z.; Xu, A.; Yang, Y.
Deposited on : 2018-09-18
Resolution : 2.80 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 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 : 4.02b-467
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 1.20.1
EDS : 3.0
Percentile statistics : 20231227.v01 (using entries in the PDB archive December 27th 2023)
CCP4 : 9.0.003 (Gargrove)
Density-Fitness : 1.0.11
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.39

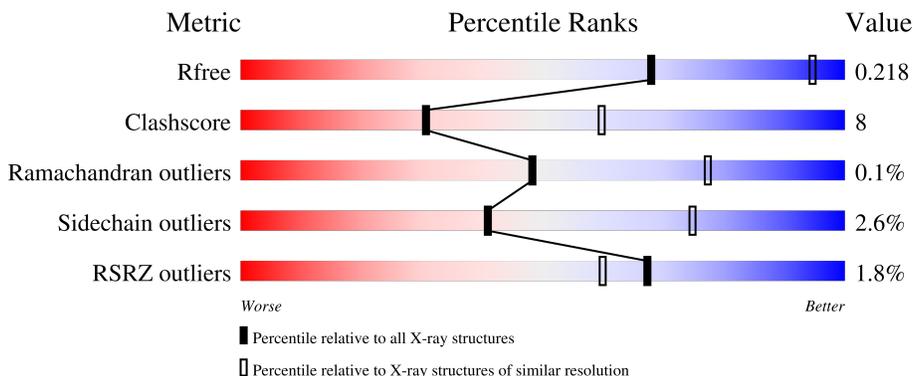
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.80 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}	164625	3657 (2.80-2.80)
Clashscore	180529	4123 (2.80-2.80)
Ramachandran outliers	177936	4071 (2.80-2.80)
Sidechain outliers	177891	4073 (2.80-2.80)
RSRZ outliers	164620	3659 (2.80-2.80)

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	235	
1	D	235	
2	B	216	
2	E	216	
3	C	144	

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Mol	Chain	Length	Quality of chain
3	F	144	
4	G	5	
5	H	4	

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
8	EPE	C	205	-	X	-	-
8	EPE	F	206	-	X	-	-

2 Entry composition [i](#)

There are 9 unique types of molecules in this entry. The entry contains 8622 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 Urelumab Fab heavy chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	D	217	Total 1623	C 1040	N 261	O 318	S 4	0	0	0
1	A	219	Total 1624	C 1040	N 261	O 319	S 4	0	0	0

- Molecule 2 is a protein called Urelumab Fab kappa chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	E	215	Total 1624	C 1018	N 274	O 328	S 4	0	0	0
2	B	215	Total 1624	C 1019	N 275	O 326	S 4	0	0	0

- Molecule 3 is a protein called Tumor necrosis factor receptor superfamily member 9.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
3	F	140	Total 975	C 579	N 176	O 199	S 21	0	0	0
3	C	141	Total 980	C 583	N 176	O 199	S 22	0	0	0

There are 12 discrepancies between the modelled and reference sequences:

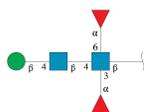
Chain	Residue	Modelled	Actual	Comment	Reference
F	163	GLU	-	expression tag	UNP Q07011
F	164	ASN	-	expression tag	UNP Q07011
F	165	LEU	-	expression tag	UNP Q07011
F	166	TYR	-	expression tag	UNP Q07011
F	167	PHE	-	expression tag	UNP Q07011
F	168	GLN	-	expression tag	UNP Q07011
C	163	GLU	-	expression tag	UNP Q07011
C	164	ASN	-	expression tag	UNP Q07011

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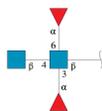
Chain	Residue	Modelled	Actual	Comment	Reference
C	165	LEU	-	expression tag	UNP Q07011
C	166	TYR	-	expression tag	UNP Q07011
C	167	PHE	-	expression tag	UNP Q07011
C	168	GLN	-	expression tag	UNP Q07011

- Molecule 4 is an oligosaccharide called beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-[alpha-L-fucopyranose-(1-3)][alpha-L-fucopyranose-(1-6)]2-acetamido-2-deoxy-beta-D-glucopyranose.



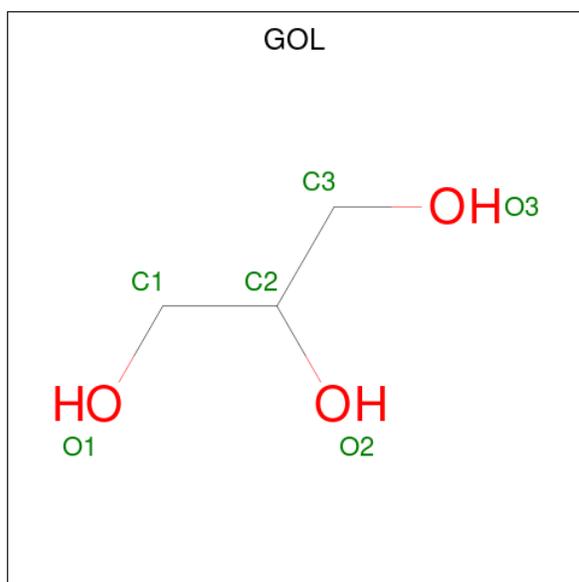
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace	
			Total	C	N				O
4	G	5	59	34	2	23	0	0	0

- Molecule 5 is an oligosaccharide called alpha-L-fucopyranose-(1-3)-[2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)][alpha-L-fucopyranose-(1-6)]2-acetamido-2-deoxy-beta-D-glucopyranose.



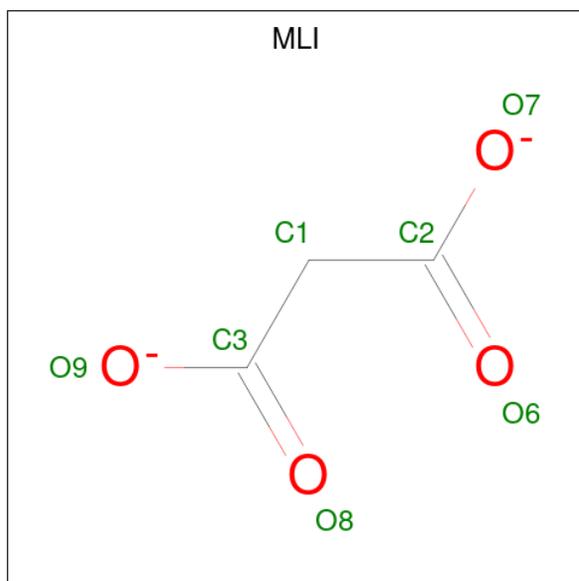
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace	
			Total	C	N				O
5	H	4	48	28	2	18	0	0	0

- Molecule 6 is GLYCEROL (three-letter code: GOL) (formula: C₃H₈O₃).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	D	1	Total	C	O	0	0
			6	3	3		

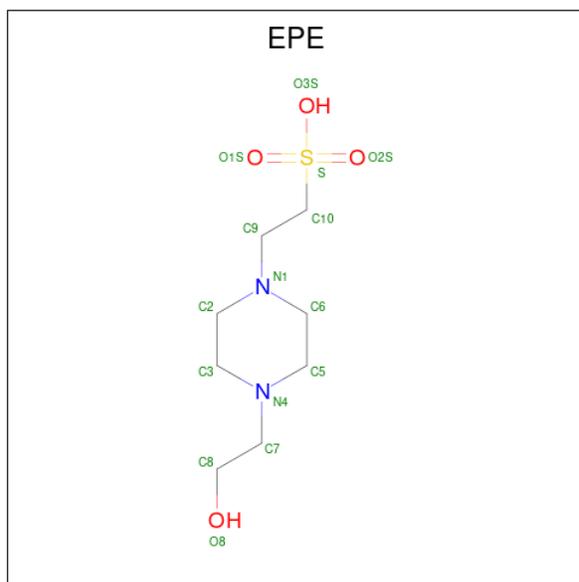
- Molecule 7 is MALONATE ION (three-letter code: MLI) (formula: $C_3H_2O_4$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
7	A	1	Total	C	O	0	0
			7	3	4		
7	B	1	Total	C	O	0	0
			7	3	4		

- Molecule 8 is 4-(2-HYDROXYETHYL)-1-PIPERAZINE ETHANESULFONIC ACID

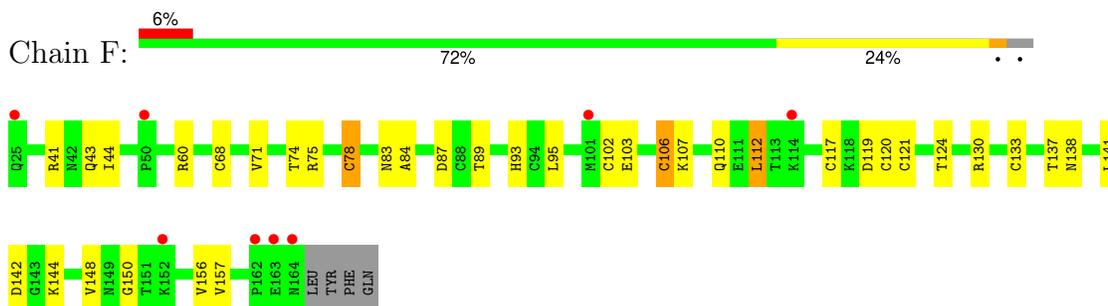
(three-letter code: EPE) (formula: C₈H₁₈N₂O₄S).



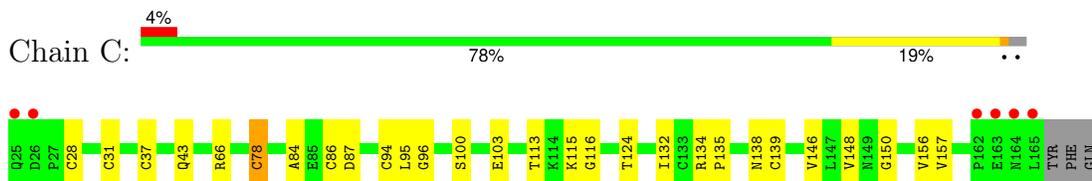
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	
			Total	C	N	O	S			
8	F	1	Total	15	8	2	4	1	0	0
8	C	1	Total	15	8	2	4	1	0	0

- Molecule 9 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
9	D	2	Total	2	0	0
9	E	5	Total	5	0	0
9	A	1	Total	1	0	0
9	B	4	Total	4	0	0
9	F	1	Total	1	0	0
9	C	2	Total	2	0	0



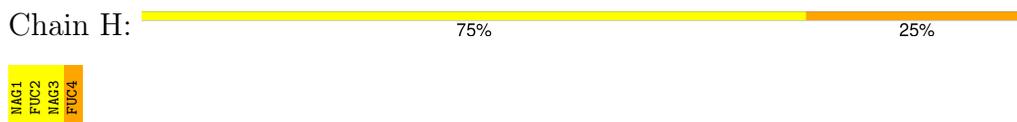
- Molecule 3: Tumor necrosis factor receptor superfamily member 9



- Molecule 4: beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-[alpha-L-fucopyranose-(1-3)][alpha-L-fucopyranose-(1-6)]2-acetamido-2-deoxy-beta-D-glucopyranose



- Molecule 5: alpha-L-fucopyranose-(1-3)-[2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)][alpha-L-fucopyranose-(1-6)]2-acetamido-2-deoxy-beta-D-glucopyranose



4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 1	Depositor
Cell constants a, b, c, α , β , γ	77.15Å 192.06Å 344.58Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	49.65 – 2.80 49.65 – 2.80	Depositor EDS
% Data completeness (in resolution range)	100.0 (49.65-2.80) 93.5 (49.65-2.80)	Depositor EDS
R_{merge}	0.22	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	0.83 (at 2.81Å)	Xtrriage
Refinement program	PHENIX 1.14_3260	Depositor
R, R_{free}	0.200 , 0.222 0.196 , 0.218	Depositor DCC
R_{free} test set	1995 reflections (3.16%)	wwPDB-VP
Wilson B-factor (Å ²)	71.7	Xtrriage
Anisotropy	0.296	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 55.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	8622	wwPDB-VP
Average B, all atoms (Å ²)	79.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.23% 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: NAG, FUC, BMA, EPE, GOL, MLI

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	0.41	0/1671	0.62	0/2296
1	D	0.49	1/1670 (0.1%)	0.65	0/2292
2	B	0.46	0/1661	0.66	0/2265
2	E	0.51	0/1661	0.66	0/2266
3	C	0.53	2/997 (0.2%)	0.63	0/1353
3	F	0.57	1/992 (0.1%)	0.82	4/1346 (0.3%)
All	All	0.49	4/8652 (0.0%)	0.67	4/11818 (0.0%)

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	F	117	CYS	CB-SG	-8.18	1.68	1.82
1	D	119	VAL	CB-CG2	-5.56	1.41	1.52
3	C	37	CYS	CB-SG	-5.37	1.73	1.81
3	C	31	CYS	CB-SG	-5.24	1.73	1.81

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	F	117	CYS	CA-CB-SG	12.02	135.63	114.00
3	F	102	CYS	CA-CB-SG	-5.46	104.17	114.00
3	F	106	CYS	CA-CB-SG	-5.43	104.22	114.00
3	F	133	CYS	CA-CB-SG	5.07	123.12	114.00

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	1624	0	1510	30	0
1	D	1623	0	1528	34	0
2	B	1624	0	1553	17	0
2	E	1624	0	1546	21	1
3	C	980	0	817	17	0
3	F	975	0	822	25	0
4	G	59	0	52	0	0
5	H	48	0	43	2	0
6	D	6	0	8	0	0
7	A	7	0	2	0	0
7	B	7	0	2	0	0
8	C	15	0	18	5	0
8	F	15	0	18	1	0
9	A	1	0	0	0	0
9	B	4	0	0	0	0
9	C	2	0	0	0	0
9	D	2	0	0	0	0
9	E	5	0	0	0	0
9	F	1	0	0	0	0
All	All	8622	0	7919	136	1

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 8.

The worst 5 of 136 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)
3:C:148:VAL:HG12	3:C:157:VAL:HG22	1.52	0.91
2:E:29:VAL:HG11	2:E:90:GLN:HG3	1.64	0.77
1:D:205:ASN:ND2	1:D:216:ASP:OD1	2.18	0.76
2:B:39:LYS:HD3	2:B:84:ALA:HB2	1.68	0.76
3:F:95:LEU:HD12	3:F:103:GLU:HB2	1.66	0.76

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:E:49:TYR:OH	2:E:49:TYR:OH[3_554]	1.66	0.54

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	215/235 (92%)	198 (92%)	16 (7%)	1 (0%)	25	56
1	D	213/235 (91%)	202 (95%)	11 (5%)	0	100	100
2	B	213/216 (99%)	207 (97%)	6 (3%)	0	100	100
2	E	213/216 (99%)	206 (97%)	7 (3%)	0	100	100
3	C	139/144 (96%)	129 (93%)	10 (7%)	0	100	100
3	F	138/144 (96%)	125 (91%)	13 (9%)	0	100	100
All	All	1131/1190 (95%)	1067 (94%)	63 (6%)	1 (0%)	48	77

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	16	GLU

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	172/199 (86%)	167 (97%)	5 (3%)	37	71
1	D	175/199 (88%)	170 (97%)	5 (3%)	37	71

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
2	B	177/186 (95%)	174 (98%)	3 (2%)	56	84
2	E	177/186 (95%)	171 (97%)	6 (3%)	32	66
3	C	103/126 (82%)	101 (98%)	2 (2%)	52	82
3	F	104/126 (82%)	101 (97%)	3 (3%)	37	71
All	All	908/1022 (89%)	884 (97%)	24 (3%)	41	75

5 of 24 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	123	SER
2	B	90	GLN
2	B	65	SER
2	B	178	SER
2	E	63	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 11 such sidechains are listed below:

Mol	Chain	Res	Type
3	F	67	GLN
3	F	110	GLN
3	C	51	ASN
3	C	39	ASN
1	A	212	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](#)

9 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
4	NAG	G	1	4,3	14,14,15	0.96	2 (14%)	17,19,21	0.67	0
4	NAG	G	2	4	14,14,15	0.51	0	17,19,21	0.69	0
4	BMA	G	3	4	11,11,12	1.35	3 (27%)	15,15,17	0.99	2 (13%)
4	FUC	G	4	4	10,10,11	1.45	2 (20%)	14,14,16	1.29	2 (14%)
4	FUC	G	5	4	10,10,11	1.15	1 (10%)	14,14,16	1.01	1 (7%)
5	NAG	H	1	3,5	14,14,15	0.83	1 (7%)	17,19,21	0.83	1 (5%)
5	FUC	H	2	5	10,10,11	1.01	1 (10%)	14,14,16	1.30	2 (14%)
5	NAG	H	3	5	14,14,15	0.44	0	17,19,21	0.83	1 (5%)
5	FUC	H	4	5	10,10,11	1.33	1 (10%)	14,14,16	0.71	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	NAG	G	1	4,3	-	2/6/23/26	0/1/1/1
4	NAG	G	2	4	-	1/6/23/26	0/1/1/1
4	BMA	G	3	4	-	2/2/19/22	0/1/1/1
4	FUC	G	4	4	-	-	0/1/1/1
4	FUC	G	5	4	-	-	0/1/1/1
5	NAG	H	1	3,5	-	0/6/23/26	0/1/1/1
5	FUC	H	2	5	-	-	0/1/1/1
5	NAG	H	3	5	-	0/6/23/26	0/1/1/1
5	FUC	H	4	5	-	-	0/1/1/1

The worst 5 of 11 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	H	4	FUC	C2-C3	3.10	1.57	1.52
5	H	1	NAG	O5-C1	-2.98	1.38	1.43
4	G	1	NAG	O5-C1	-2.68	1.39	1.43
4	G	4	FUC	C2-C3	2.54	1.56	1.52
4	G	3	BMA	C1-C2	2.38	1.57	1.52

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

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	H	2	FUC	C1-O5-C5	2.84	119.67	112.97
5	H	2	FUC	C1-C2-C3	2.82	113.76	109.64
4	G	4	FUC	C1-C2-C3	2.79	113.71	109.64
4	G	4	FUC	C1-O5-C5	2.66	119.25	112.97
5	H	1	NAG	O3-C3-C2	2.54	114.67	109.40

There are no chirality outliers.

All (5) torsion outliers are listed below:

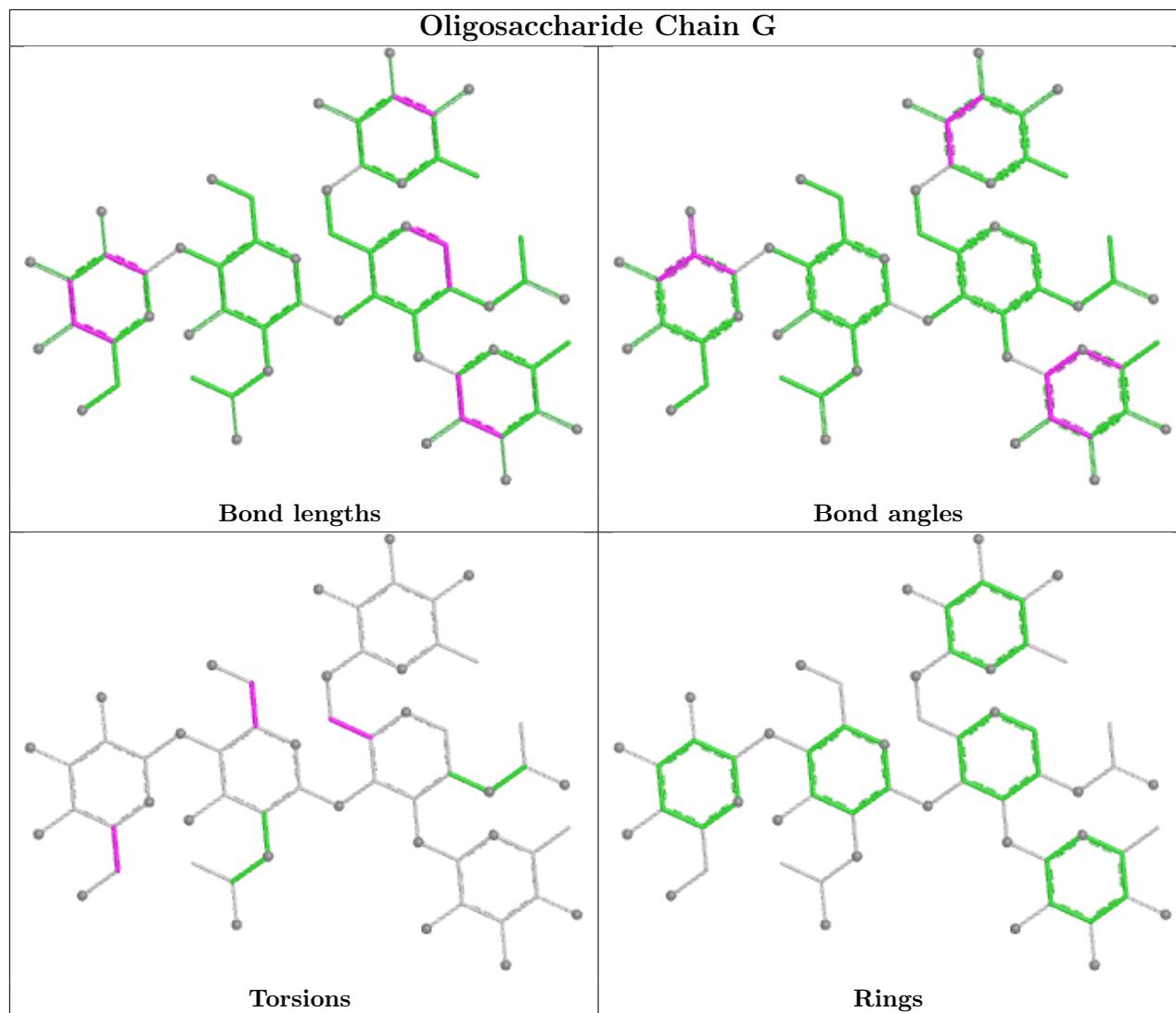
Mol	Chain	Res	Type	Atoms
4	G	3	BMA	O5-C5-C6-O6
4	G	3	BMA	C4-C5-C6-O6
4	G	1	NAG	O5-C5-C6-O6
4	G	1	NAG	C4-C5-C6-O6
4	G	2	NAG	O5-C5-C6-O6

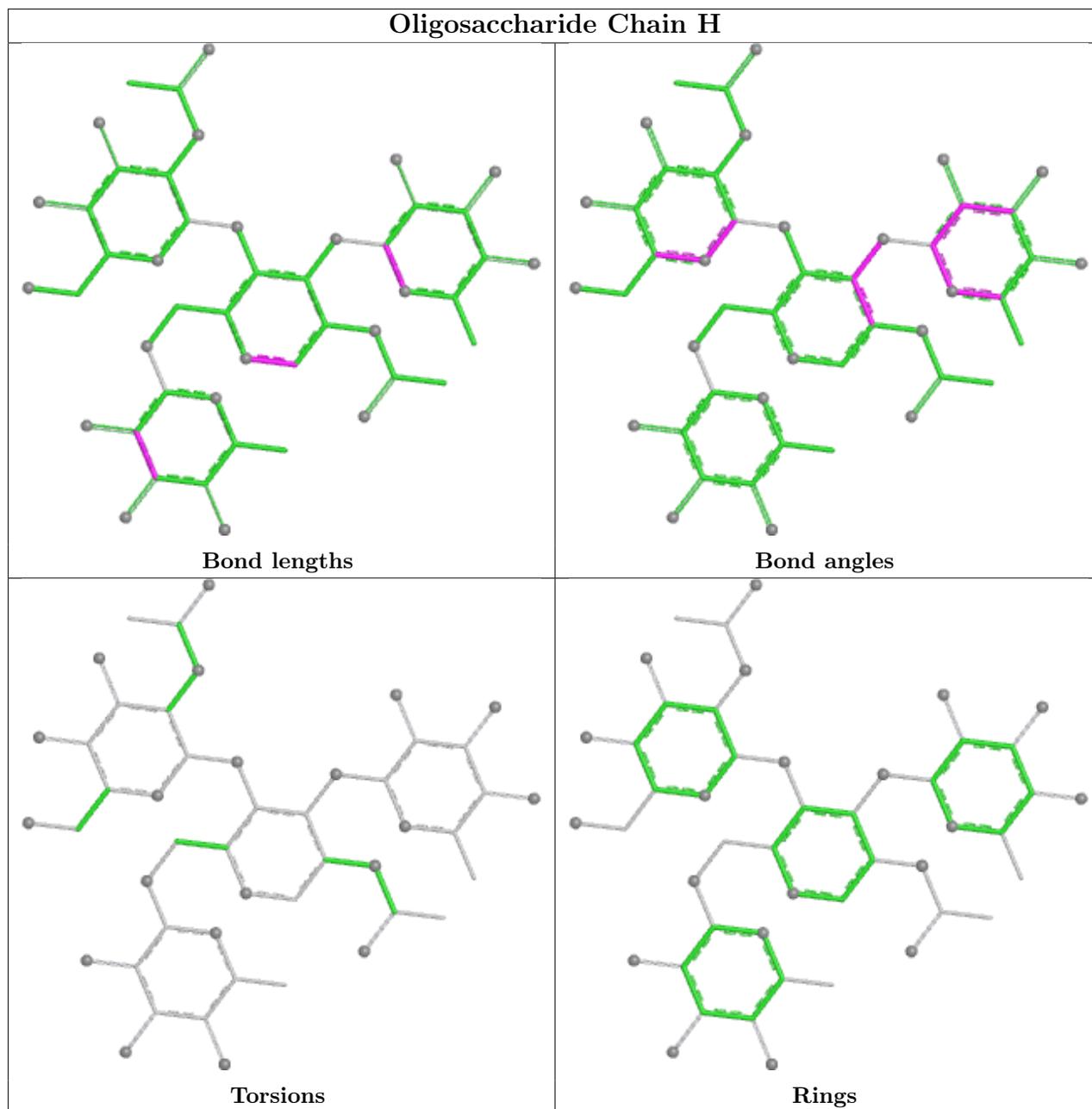
There are no ring outliers.

1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	H	4	FUC	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 oligosaccharide.





5.6 Ligand geometry [i](#)

5 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
8	EPE	F	206	-	15,15,15	3.71	14 (93%)	19,20,20	3.17	12 (63%)
7	MLI	B	301	-	6,6,6	1.61	1 (16%)	7,7,7	1.36	1 (14%)
6	GOL	D	301	-	5,5,5	0.77	0	5,5,5	0.91	0
8	EPE	C	205	-	15,15,15	4.41	11 (73%)	19,20,20	3.10	10 (52%)
7	MLI	A	301	-	6,6,6	1.45	0	7,7,7	1.39	1 (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
8	EPE	F	206	-	-	1/9/19/19	0/1/1/1
7	MLI	B	301	-	-	0/4/4/4	-
6	GOL	D	301	-	-	2/4/4/4	-
8	EPE	C	205	-	-	4/9/19/19	0/1/1/1
7	MLI	A	301	-	-	2/4/4/4	-

The worst 5 of 26 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
8	C	205	EPE	O2S-S	8.18	1.68	1.45
8	C	205	EPE	C10-S	7.59	1.88	1.77
8	C	205	EPE	O1S-S	6.89	1.64	1.45
8	F	206	EPE	O2S-S	6.19	1.62	1.45
8	F	206	EPE	O1S-S	5.82	1.61	1.45

The worst 5 of 24 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
8	C	205	EPE	O3S-S-C10	7.33	120.35	106.00
8	C	205	EPE	O2S-S-C10	-5.85	97.89	106.73
8	F	206	EPE	C3-C2-N1	-5.58	99.39	110.65
8	F	206	EPE	O2S-S-C10	-5.06	99.08	106.73
8	F	206	EPE	C6-C5-N4	-4.98	100.61	110.65

There are no chirality outliers.

5 of 9 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
6	D	301	GOL	O1-C1-C2-O2
6	D	301	GOL	O1-C1-C2-C3
8	C	205	EPE	C10-C9-N1-C6
8	C	205	EPE	C8-C7-N4-C5
8	F	206	EPE	C8-C7-N4-C5

There are no ring outliers.

2 monomers are involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
8	F	206	EPE	1	0
8	C	205	EPE	5	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 [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	219/235 (93%)	0.15	3 (1%) 73 66	58, 88, 108, 126	0
1	D	217/235 (92%)	-0.31	3 (1%) 73 66	48, 70, 103, 125	0
2	B	215/216 (99%)	-0.45	0 100 100	51, 68, 90, 108	0
2	E	215/216 (99%)	-0.56	1 (0%) 87 83	46, 64, 93, 126	0
3	C	141/144 (97%)	0.38	6 (4%) 40 32	75, 95, 114, 130	0
3	F	140/144 (97%)	0.52	8 (5%) 30 24	55, 96, 130, 145	0
All	All	1147/1190 (96%)	-0.11	21 (1%) 67 60	46, 78, 110, 145	0

The worst 5 of 21 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
3	F	164	ASN	7.4
3	C	165	LEU	6.7
1	A	135	SER	4.7
3	F	162	PRO	4.5
3	F	163	GLU	3.8

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

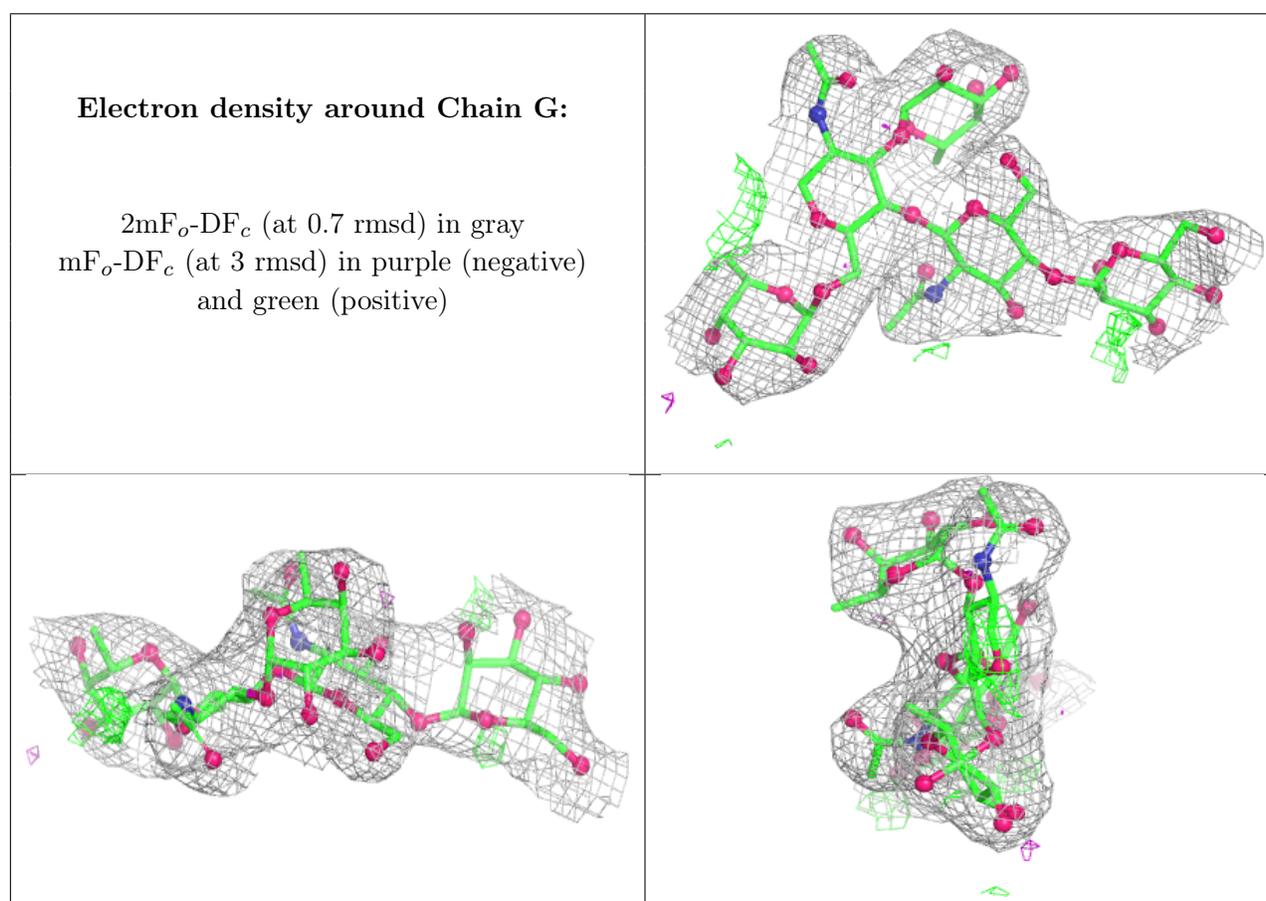
There are no non-standard protein/DNA/RNA residues in this entry.

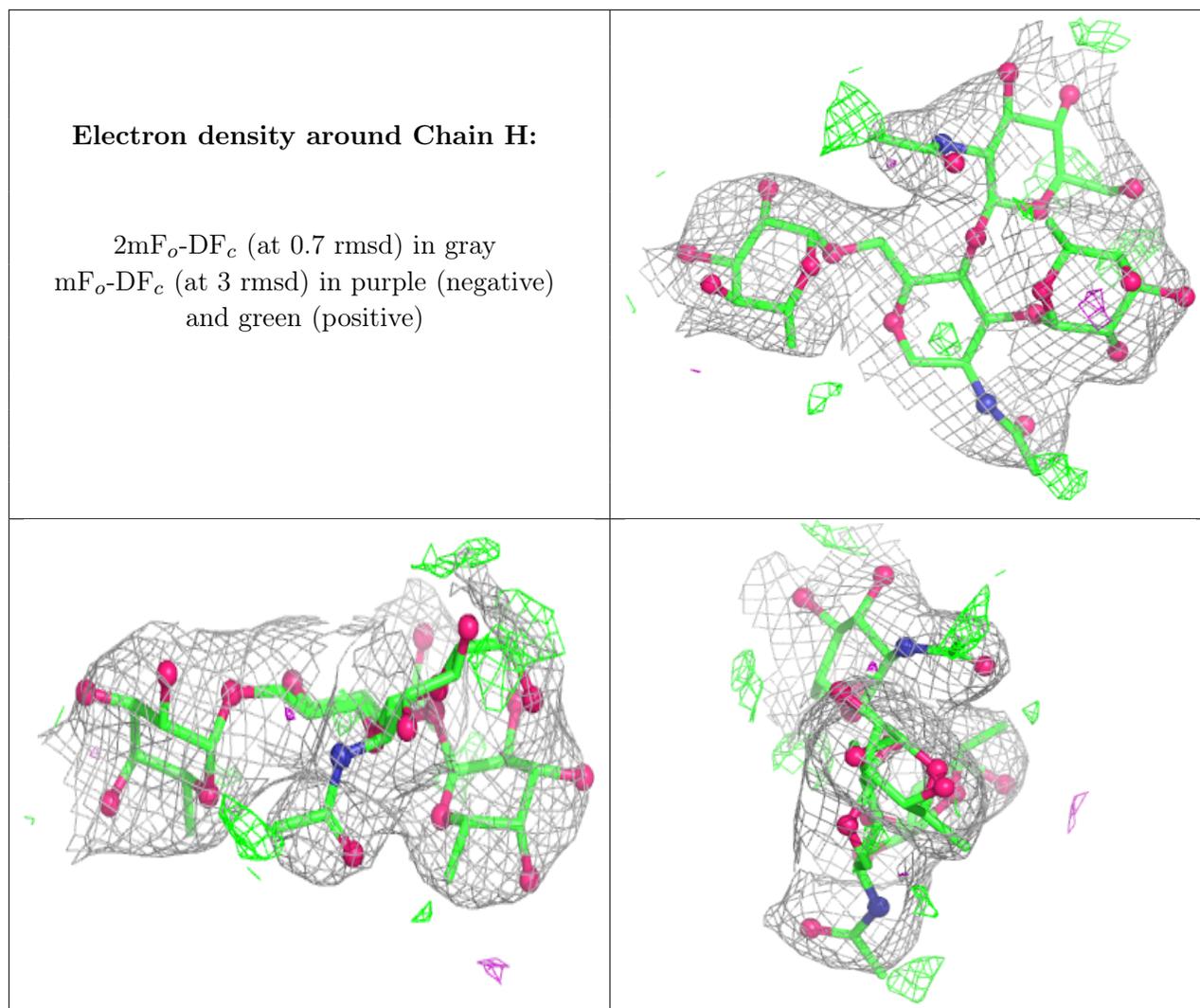
6.3 Carbohydrates [i](#)

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
4	BMA	G	3	11/12	0.58	0.14	127,132,140,142	0
5	FUC	H	2	10/11	0.75	0.16	126,139,142,144	0
5	NAG	H	3	14/15	0.75	0.16	110,131,138,138	0
5	NAG	H	1	14/15	0.90	0.09	106,112,124,128	0
5	FUC	H	4	10/11	0.92	0.14	91,109,115,118	0
4	NAG	G	2	14/15	0.93	0.09	75,104,113,125	0
4	FUC	G	5	10/11	0.93	0.11	73,85,92,98	0
4	FUC	G	4	10/11	0.94	0.10	93,101,105,105	0
4	NAG	G	1	14/15	0.96	0.07	85,91,102,104	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(Å ²)	Q<0.9
7	MLI	B	301	7/7	0.73	0.12	103,108,114,124	0
7	MLI	A	301	7/7	0.75	0.12	110,127,129,131	0
6	GOL	D	301	6/6	0.83	0.14	86,92,95,95	0
8	EPE	C	205	15/15	0.88	0.18	92,96,108,114	0
8	EPE	F	206	15/15	0.93	0.15	71,81,99,102	0

6.5 Other polymers [i](#)

There are no such residues in this entry.