



wwPDB X-ray Structure Validation Summary Report ⓘ

Mar 10, 2026 – 11:02 AM UTC

PDB ID : 4MWF / pdb_00004mwf
Title : Structure of Hepatitis C Virus Envelope Glycoprotein E2 core bound to broadly neutralizing antibody AR3C
Authors : Kong, L.; Wilson, I.A.; Law, M.
Deposited on : 2013-09-24
Resolution : 2.65 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	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

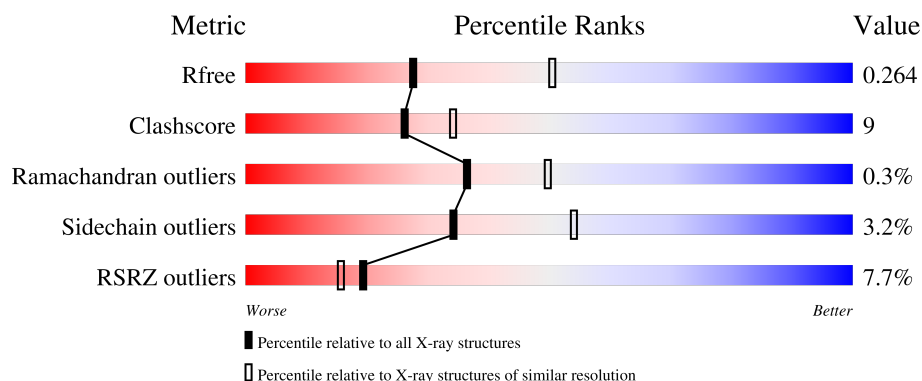
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.65 Å.

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



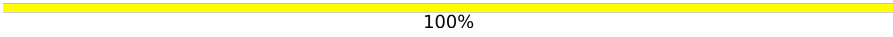
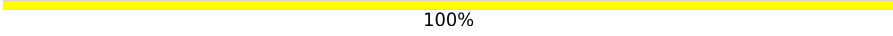
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	2053 (2.66-2.62)
Clashscore	190562	2097 (2.66-2.62)
Ramachandran outliers	187476	2066 (2.66-2.62)
Sidechain outliers	187428	2066 (2.66-2.62)
RSRZ outliers	180081	2052 (2.66-2.62)

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	233	11% 77% 17% • •
1	H	233	6% 83% 14% • •
2	B	214	3% 86% 12% •
2	L	214	2% 86% 13% •
3	C	212	8% 45% 30% • 20%

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Mol	Chain	Length	Quality of chain
3	D	212	
4	E	8	
5	F	2	
5	G	2	

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 9433 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 Fab AR3C heavy chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	223	Total	C	N	O	S	0	0	0
			1663	1047	282	327	7			
1	H	228	Total	C	N	O	S	0	0	0
			1697	1065	288	337	7			

- Molecule 2 is a protein called Fab AR3C light chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	211	Total	C	N	O	S	0	0	0
			1612	1005	277	326	4			
2	L	212	Total	C	N	O	S	0	0	0
			1620	1011	278	327	4			

- Molecule 3 is a protein called Envelope glycoprotein E2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	169	Total	C	N	O	S	0	0	0
			1288	818	218	235	17			
3	D	171	Total	C	N	O	S	0	0	0
			1305	831	222	235	17			

There are 16 discrepancies between the modelled and reference sequences:

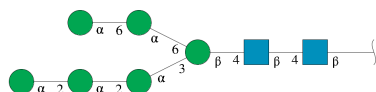
Chain	Residue	Modelled	Actual	Comment	Reference
C	448	ASP	ASN	engineered mutation	UNP P27958
C	482	GLY	-	linker	UNP P27958
C	483	SER	-	linker	UNP P27958
C	484	SER	-	linker	UNP P27958
C	485	GLY	-	linker	UNP P27958
C	576	ASP	ASN	engineered mutation	UNP P27958
C	589	HIS	TYR	variant	UNP P27958
C	602	TRP	ARG	variant	UNP P27958

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Chain	Residue	Modelled	Actual	Comment	Reference
D	448	ASP	ASN	engineered mutation	UNP P27958
D	482	GLY	-	linker	UNP P27958
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D	485	GLY	-	linker	UNP P27958
D	576	ASP	ASN	engineered mutation	UNP P27958
D	589	HIS	TYR	variant	UNP P27958
D	602	TRP	ARG	variant	UNP P27958

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



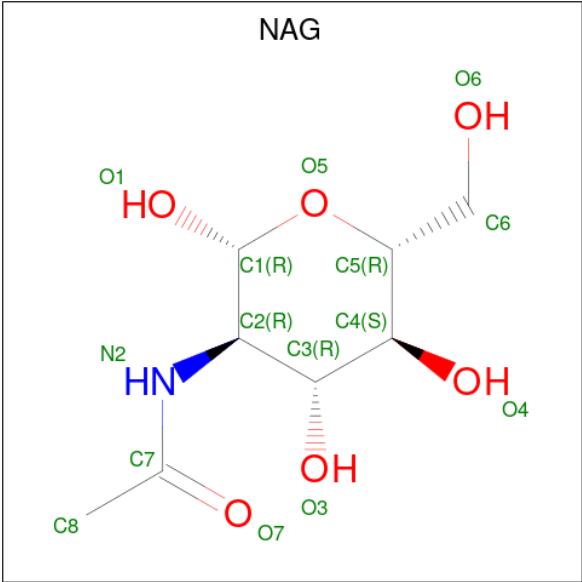
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
4	E	8	Total	C	N	O	0	0	0
			94	52	2	40			

- Molecule 5 is an oligosaccharide called 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
5	F	2	Total	C	N	O	0	0	0
			28	16	2	10			
5	G	2	Total	C	N	O	0	0	0
			28	16	2	10			

- Molecule 6 is 2-acetamido-2-deoxy-beta-D-glucopyranose (CCD ID: NAG) (formula: C₈H₁₅NO₆).

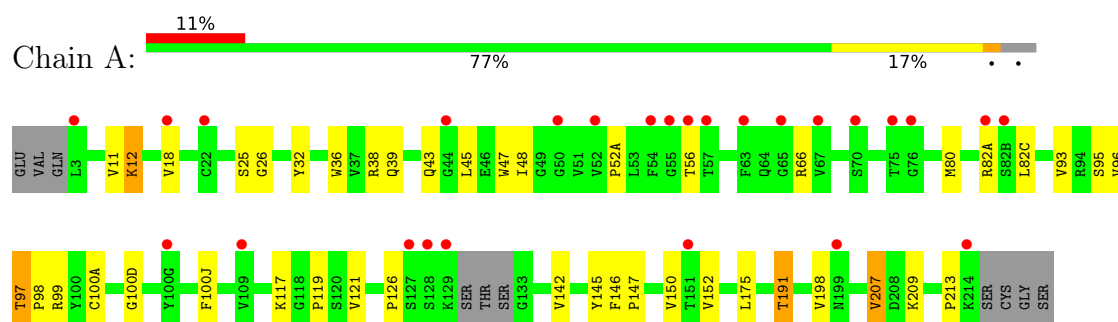


Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
6	C	1	Total	C	N	O	0	0
			14	8	1	5		
6	C	1	Total	C	N	O	0	0
			14	8	1	5		
6	C	1	Total	C	N	O	0	0
			14	8	1	5		
6	D	1	Total	C	N	O	0	0
			14	8	1	5		
6	D	1	Total	C	N	O	0	0
			14	8	1	5		
6	D	1	Total	C	N	O	0	0
			14	8	1	5		

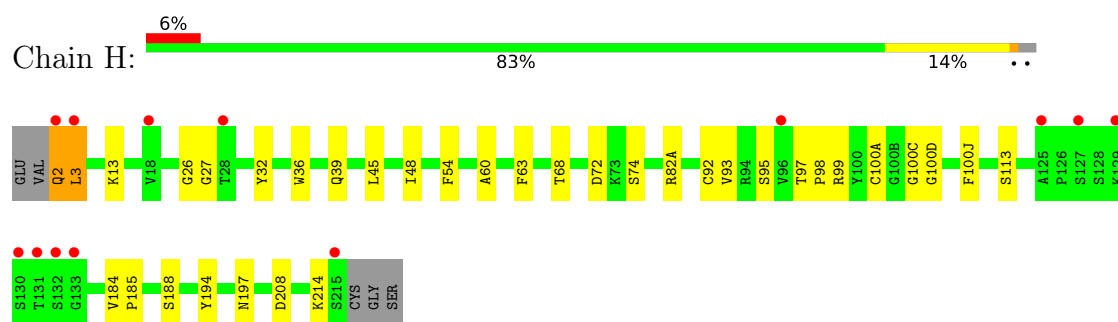
3 Residue-property plots [i](#)

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.

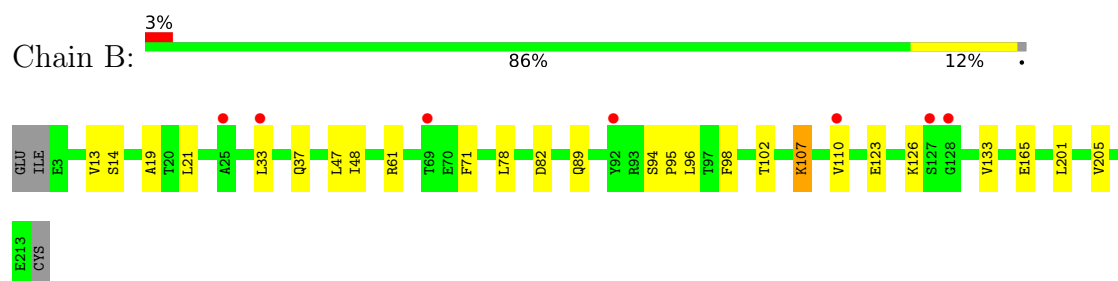
- Molecule 1: Fab AR3C heavy chain



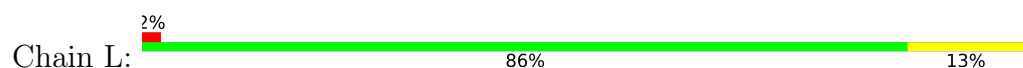
- Molecule 1: Fab AR3C heavy chain

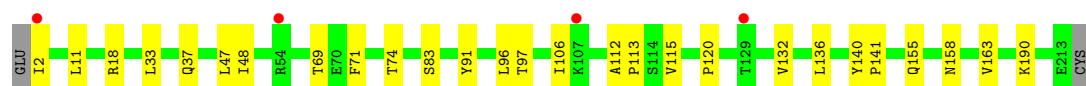


- Molecule 2: Fab AR3C light chain

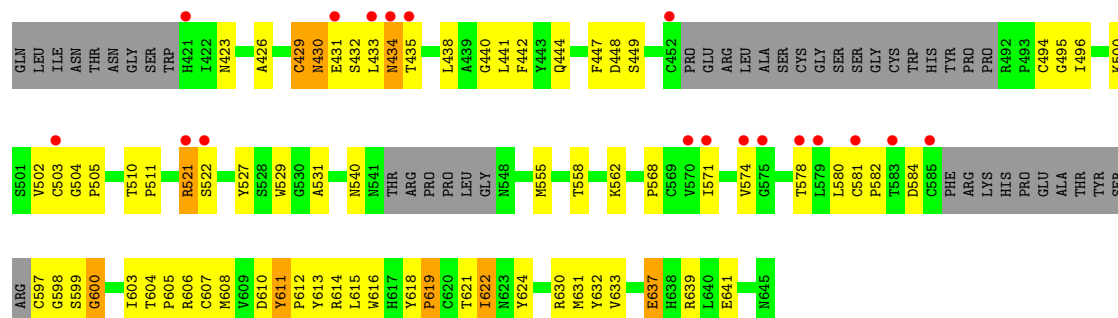


- Molecule 2: Fab AR3C light chain

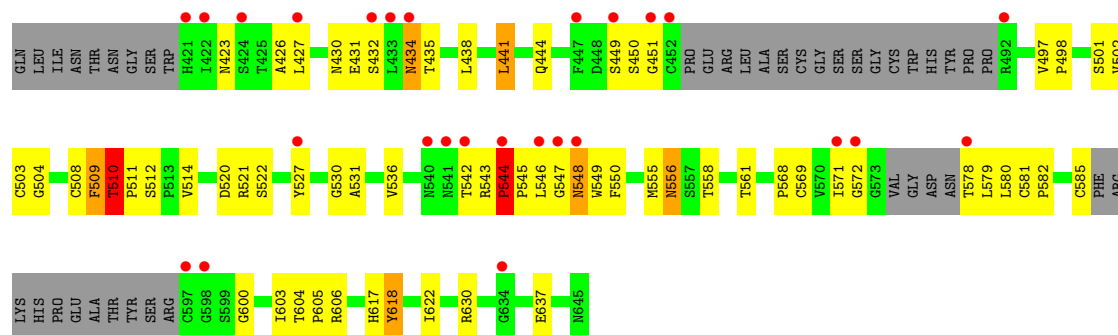




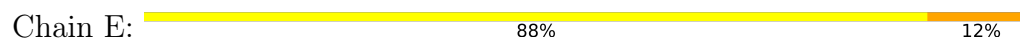
• Molecule 3: Envelope glycoprotein E2



• Molecule 3: Envelope glycoprotein E2



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



• Molecule 5: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose



- Molecule 5: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain G:

100%

MAG1
MAG2

4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	47.13Å 166.55Å 209.96Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	49.08 – 2.65 49.08 – 2.65	Depositor EDS
% Data completeness (in resolution range)	92.7 (49.08-2.65) 92.8 (49.08-2.65)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.77 (at 2.65Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.8.2_1309)	Depositor
R, R_{free}	0.231 , 0.270 0.231 , 0.264	Depositor DCC
R_{free} test set	2323 reflections (4.69%)	wwPDB-VP
Wilson B-factor (Å ²)	55.5	Xtriage
Anisotropy	0.761	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 47.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	9433	wwPDB-VP
Average B, all atoms (Å ²)	71.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 10.00% 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: BMA, MAN, NAG

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	0.42	0/1702	0.83	1/2314 (0.0%)
1	H	0.48	0/1737	0.87	1/2363 (0.0%)
2	B	0.40	0/1646	0.82	1/2235 (0.0%)
2	L	0.48	0/1654	0.85	1/2246 (0.0%)
3	C	0.56	2/1328 (0.2%)	1.02	8/1816 (0.4%)
3	D	0.48	0/1347	1.14	13/1843 (0.7%)
All	All	0.47	2/9414 (0.0%)	0.92	25/12817 (0.2%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	C	619	PRO	N-CD	5.25	1.55	1.47
3	C	612	PRO	N-CD	5.21	1.55	1.47

The worst 5 of 25 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	600	GLY	N-CA-C	7.96	121.56	112.00
3	D	509	PHE	N-CA-C	7.01	120.35	109.07
3	C	430	ASN	N-CA-C	6.69	119.49	108.99
3	D	543	ARG	CA-C-N	6.68	127.27	120.38
3	D	543	ARG	C-N-CA	6.68	127.27	120.38

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	1663	0	1626	32	0
1	H	1697	0	1653	25	0
2	B	1612	0	1570	12	0
2	L	1620	0	1581	11	0
3	C	1288	0	1193	56	0
3	D	1305	0	1219	46	0
4	E	94	0	79	1	0
5	F	28	0	25	0	0
5	G	28	0	25	5	0
6	C	42	0	39	1	0
6	D	56	0	52	0	0
All	All	9433	0	9062	173	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 9.

The worst 5 of 173 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:433:LEU:HD12	1:H:99:ARG:HD2	1.43	1.01
1:H:97:THR:HG22	1:H:99:ARG:HH12	1.35	0.90
3:C:433:LEU:CD1	1:H:99:ARG:HD2	2.02	0.90
3:C:610:ASP:HB3	3:C:641:GLU:HB3	1.65	0.78
1:A:66:ARG:HE	1:A:82(A):ARG:HB2	1.52	0.74

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	219/233 (94%)	216 (99%)	2 (1%)	1 (0%)	24	36
1	H	226/233 (97%)	221 (98%)	5 (2%)	0	100	100
2	B	209/214 (98%)	201 (96%)	8 (4%)	0	100	100
2	L	210/214 (98%)	206 (98%)	4 (2%)	0	100	100
3	C	161/212 (76%)	152 (94%)	8 (5%)	1 (1%)	21	31
3	D	163/212 (77%)	150 (92%)	11 (7%)	2 (1%)	10	15
All	All	1188/1318 (90%)	1146 (96%)	38 (3%)	4 (0%)	36	50

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	97	THR
3	D	510	THR
3	C	510	THR
3	D	545	PRO

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	186/195 (95%)	181 (97%)	5 (3%)	39	60
1	H	191/195 (98%)	187 (98%)	4 (2%)	47	67
2	B	181/184 (98%)	176 (97%)	5 (3%)	38	58
2	L	182/184 (99%)	176 (97%)	6 (3%)	33	53
3	C	143/180 (79%)	136 (95%)	7 (5%)	22	37
3	D	145/180 (81%)	139 (96%)	6 (4%)	27	44
All	All	1028/1118 (92%)	995 (97%)	33 (3%)	34	54

5 of 33 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	L	48	ILE
2	L	69	THR
2	L	190	LYS
3	C	521	ARG
3	C	444	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 12 such sidechains are listed below:

Mol	Chain	Res	Type
3	C	541	ASN
3	D	428	ASN
1	H	164	HIS
1	H	59	ASN
2	B	137	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 ⓘ

12 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	E	1	4,3	14,14,15	0.43	0	17,19,21	0.98	1 (5%)
4	NAG	E	2	4	14,14,15	0.51	0	17,19,21	1.19	1 (5%)
4	BMA	E	3	4	11,11,12	0.84	1 (9%)	15,15,17	1.52	2 (13%)
4	MAN	E	4	4	11,11,12	0.38	0	15,15,17	1.12	2 (13%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
4	MAN	E	5	4	11,11,12	0.33	0	15,15,17	0.87	1 (6%)
4	MAN	E	6	4	11,11,12	0.20	0	15,15,17	0.73	0
4	MAN	E	7	4	11,11,12	0.88	0	15,15,17	2.26	2 (13%)
4	MAN	E	8	4	11,11,12	0.27	0	15,15,17	0.89	1 (6%)
5	NAG	F	1	3,5	14,14,15	0.51	0	17,19,21	1.10	1 (5%)
5	NAG	F	2	5	14,14,15	0.58	0	17,19,21	0.81	1 (5%)
5	NAG	G	1	3,5	14,14,15	0.28	0	17,19,21	0.57	0
5	NAG	G	2	5	14,14,15	0.30	0	17,19,21	0.55	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	E	1	4,3	-	2/6/23/26	0/1/1/1
4	NAG	E	2	4	-	2/6/23/26	0/1/1/1
4	BMA	E	3	4	-	0/2/19/22	0/1/1/1
4	MAN	E	4	4	-	2/2/19/22	0/1/1/1
4	MAN	E	5	4	-	0/2/19/22	0/1/1/1
4	MAN	E	6	4	-	2/2/19/22	0/1/1/1
4	MAN	E	7	4	-	0/2/19/22	0/1/1/1
4	MAN	E	8	4	-	2/2/19/22	0/1/1/1
5	NAG	F	1	3,5	-	3/6/23/26	0/1/1/1
5	NAG	F	2	5	-	3/6/23/26	0/1/1/1
5	NAG	G	1	3,5	-	2/6/23/26	0/1/1/1
5	NAG	G	2	5	-	2/6/23/26	0/1/1/1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	E	3	BMA	C4-C5	2.13	1.57	1.53

The worst 5 of 12 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	E	7	MAN	C1-O5-C5	7.83	122.68	112.19
4	E	3	BMA	C1-O5-C5	3.82	117.30	112.19
4	E	3	BMA	C1-C2-C3	3.48	114.71	109.64

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	F	1	NAG	O5-C1-C2	-2.94	106.74	111.29
4	E	1	NAG	C2-N2-C7	-2.85	119.07	122.90

There are no chirality outliers.

5 of 20 torsion outliers are listed below:

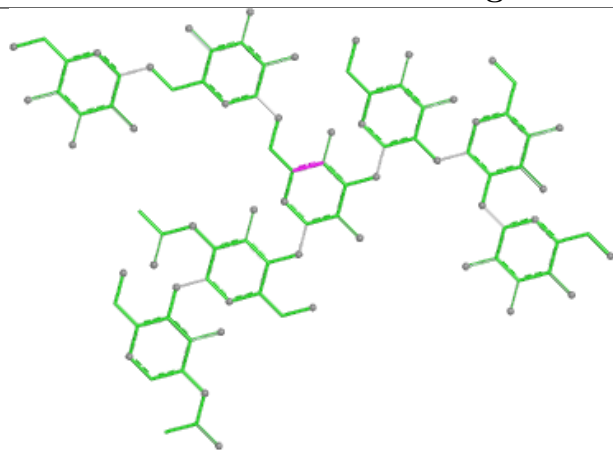
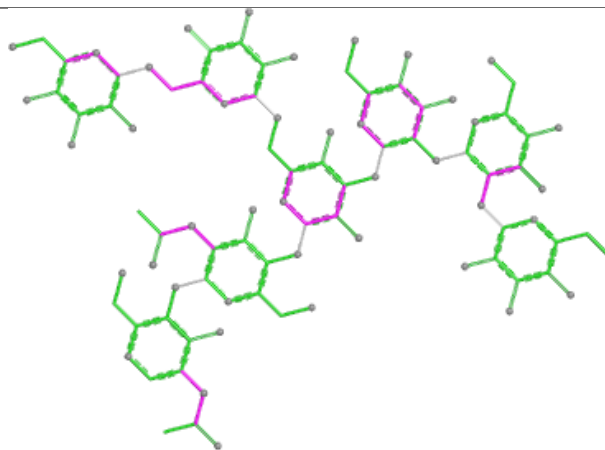
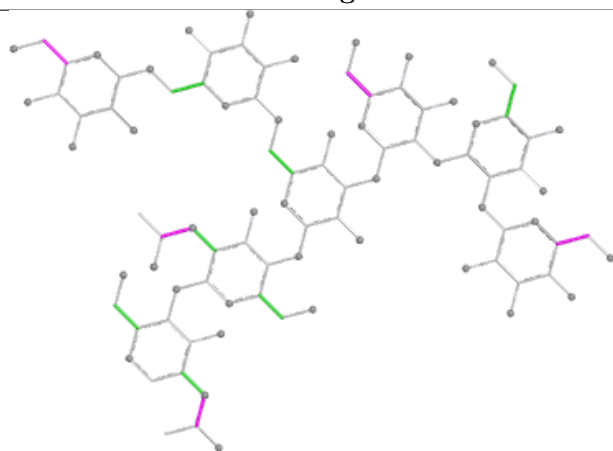
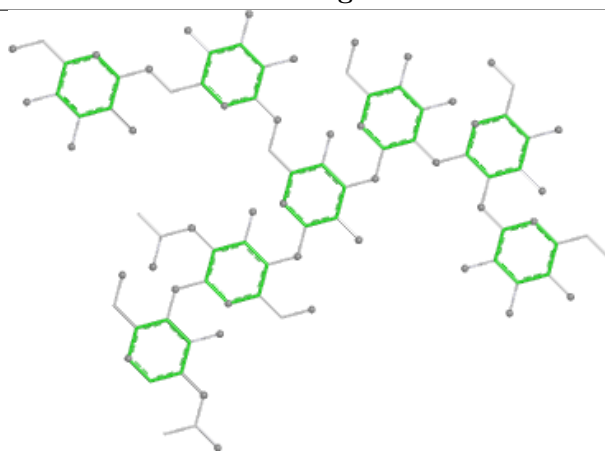
Mol	Chain	Res	Type	Atoms
4	E	4	MAN	O5-C5-C6-O6
5	G	1	NAG	O5-C5-C6-O6
5	G	2	NAG	O5-C5-C6-O6
5	F	2	NAG	C8-C7-N2-C2
5	F	2	NAG	O7-C7-N2-C2

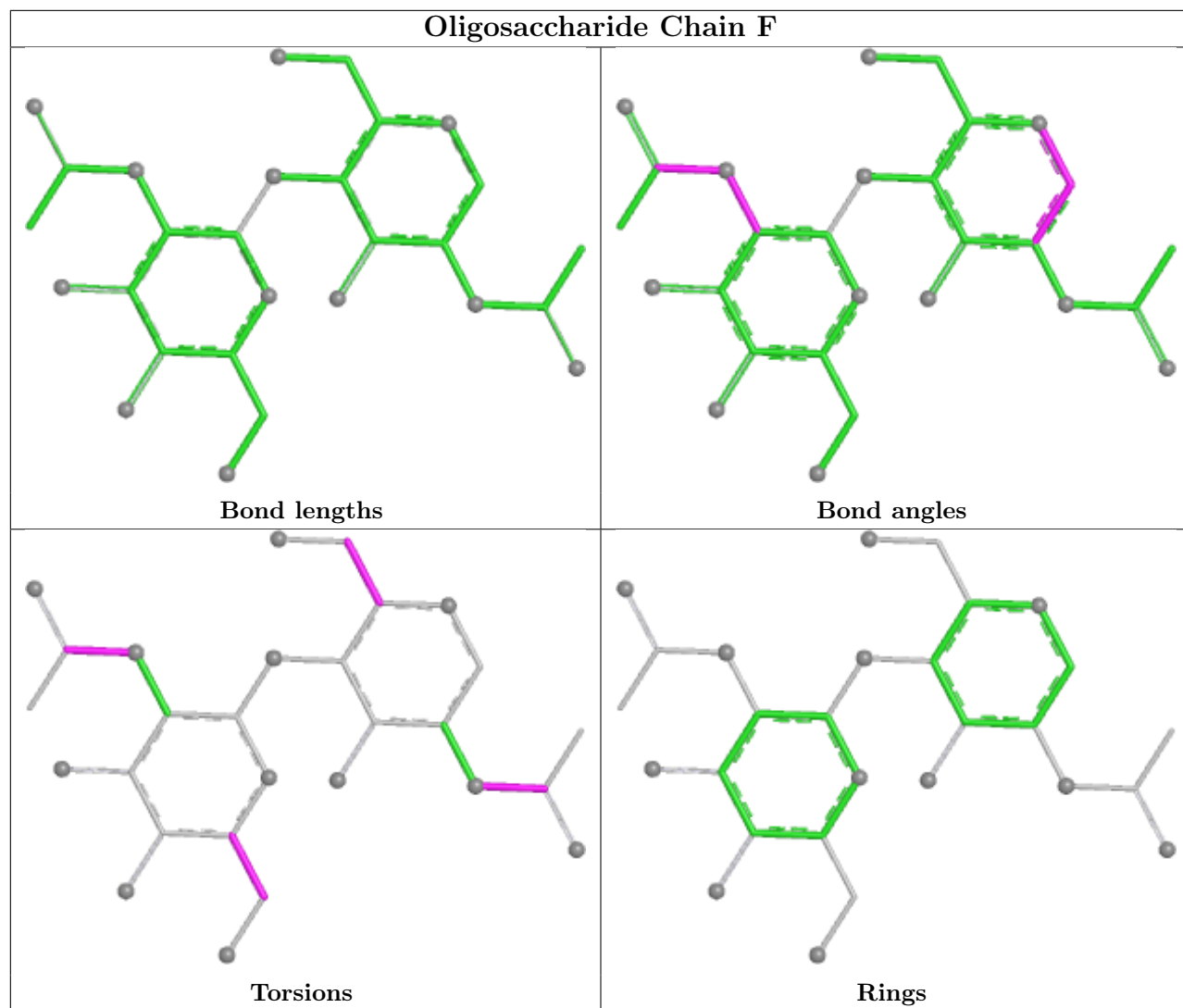
There are no ring outliers.

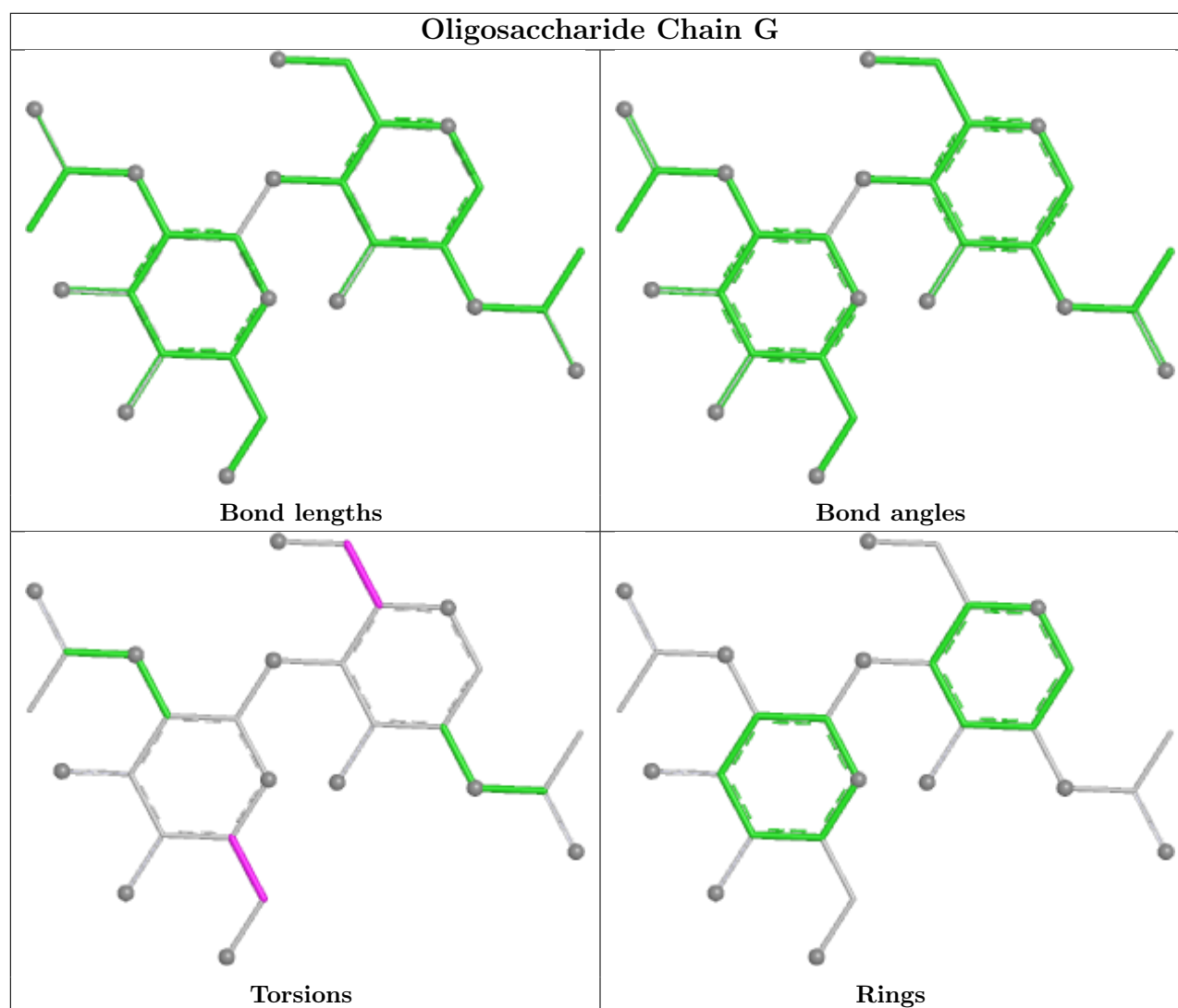
4 monomers are involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	G	1	NAG	5	0
5	G	2	NAG	4	0
4	E	5	MAN	1	0
4	E	6	MAN	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.

Oligosaccharide Chain E**Bond lengths****Bond angles****Torsions****Rings**





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	$\# Z > 2$
6	NAG	D	703	3	14,14,15	0.30	0	17,19,21	0.56	0
6	NAG	C	713	3	14,14,15	0.46	0	17,19,21	1.02	1 (5%)
6	NAG	D	706	3	14,14,15	0.55	0	17,19,21	0.77	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
6	NAG	C	712	3	14,14,15	0.28	0	17,19,21	0.53	0
6	NAG	C	711	3	14,14,15	0.50	0	17,19,21	0.85	1 (5%)
6	NAG	D	701	3	14,14,15	0.49	0	17,19,21	0.91	1 (5%)
6	NAG	D	702	3	14,14,15	0.72	0	17,19,21	0.90	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
6	NAG	D	703	3	-	3/6/23/26	0/1/1/1
6	NAG	C	713	3	-	0/6/23/26	0/1/1/1
6	NAG	D	706	3	-	2/6/23/26	0/1/1/1
6	NAG	C	712	3	-	2/6/23/26	0/1/1/1
6	NAG	C	711	3	-	0/6/23/26	0/1/1/1
6	NAG	D	701	3	-	2/6/23/26	0/1/1/1
6	NAG	D	702	3	-	1/6/23/26	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(°)
6	C	713	NAG	C1-O5-C5	3.19	116.46	112.19
6	D	701	NAG	C1-O5-C5	2.44	115.46	112.19
6	C	711	NAG	O5-C5-C6	2.24	112.02	107.66

There are no chirality outliers.

5 of 10 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
6	C	712	NAG	O7-C7-N2-C2
6	C	712	NAG	C8-C7-N2-C2
6	D	701	NAG	O5-C5-C6-O6
6	D	703	NAG	C8-C7-N2-C2
6	D	703	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
6	C	712	NAG	1	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data [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	223/233 (95%)	0.89	26 (11%) 9 7	46, 77, 109, 123	0
1	H	228/233 (97%)	0.39	13 (5%) 29 24	39, 55, 96, 134	0
2	B	211/214 (98%)	0.52	7 (3%) 49 43	48, 72, 95, 114	0
2	L	212/214 (99%)	0.34	4 (1%) 66 62	31, 57, 76, 101	1 (0%)
3	C	169/212 (79%)	0.75	18 (10%) 11 9	45, 69, 123, 165	0
3	D	171/212 (80%)	1.02	26 (15%) 5 4	50, 81, 120, 141	0
All	All	1214/1318 (92%)	0.64	94 (7%) 19 16	31, 66, 109, 165	1 (0%)

The worst 5 of 94 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	L	54	ARG	5.5
1	H	131	THR	4.7
1	H	3	LEU	4.7
3	D	547	GLY	4.4
1	A	55	GLY	4.3

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

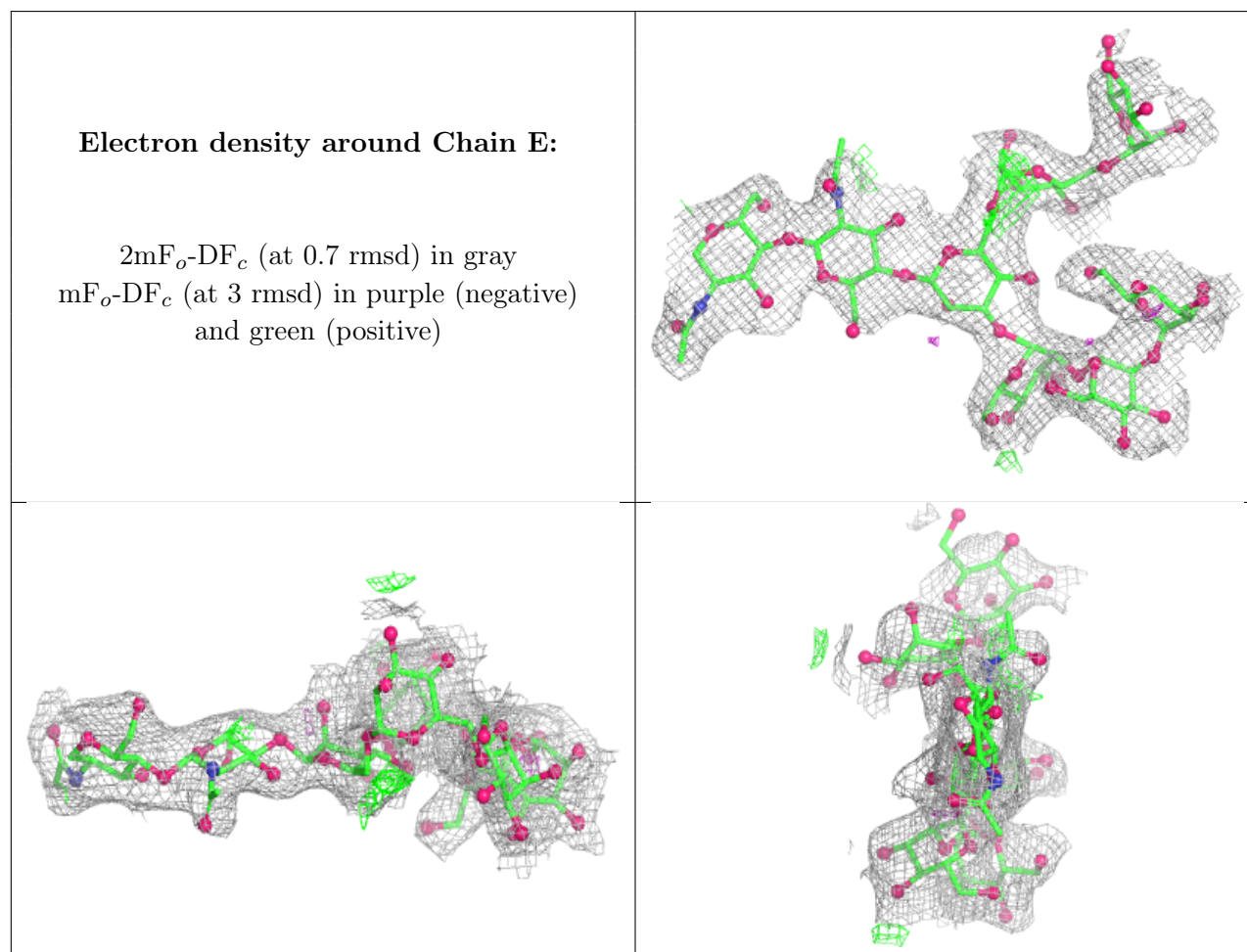
There are no non-standard protein/DNA/RNA residues in this entry.

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.

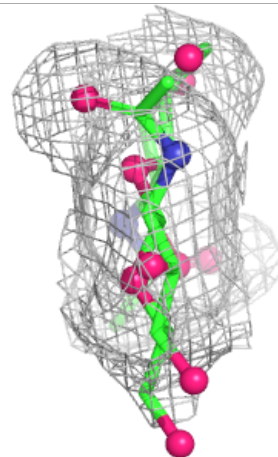
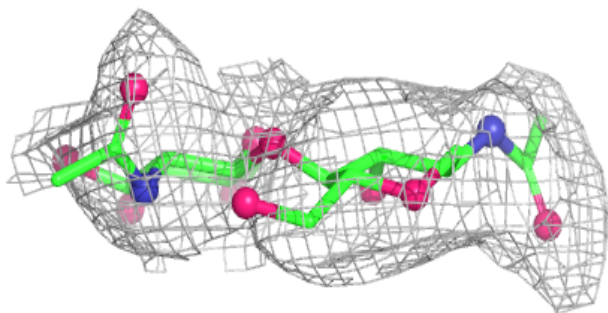
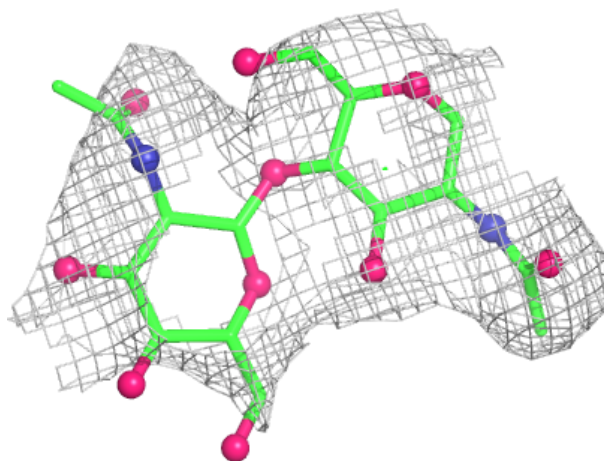
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
4	NAG	E	1	14/15	-	-	60,71,86,92	0
4	NAG	E	2	14/15	-	-	78,86,102,108	0
4	BMA	E	3	11/12	-	-	78,86,99,107	0
4	MAN	E	4	11/12	-	-	55,69,80,90	0
4	MAN	E	5	11/12	-	-	64,78,85,91	0
4	MAN	E	6	11/12	-	-	81,98,105,108	0
4	MAN	E	7	11/12	-	-	96,111,124,130	0
4	MAN	E	8	11/12	-	-	131,135,140,148	0
5	NAG	F	1	14/15	-	-	63,88,114,115	0
5	NAG	F	2	14/15	-	-	102,121,136,144	0
5	NAG	G	2	14/15	0.82	0.14	113,137,153,153	0
5	NAG	G	1	14/15	0.91	0.10	63,93,112,114	0

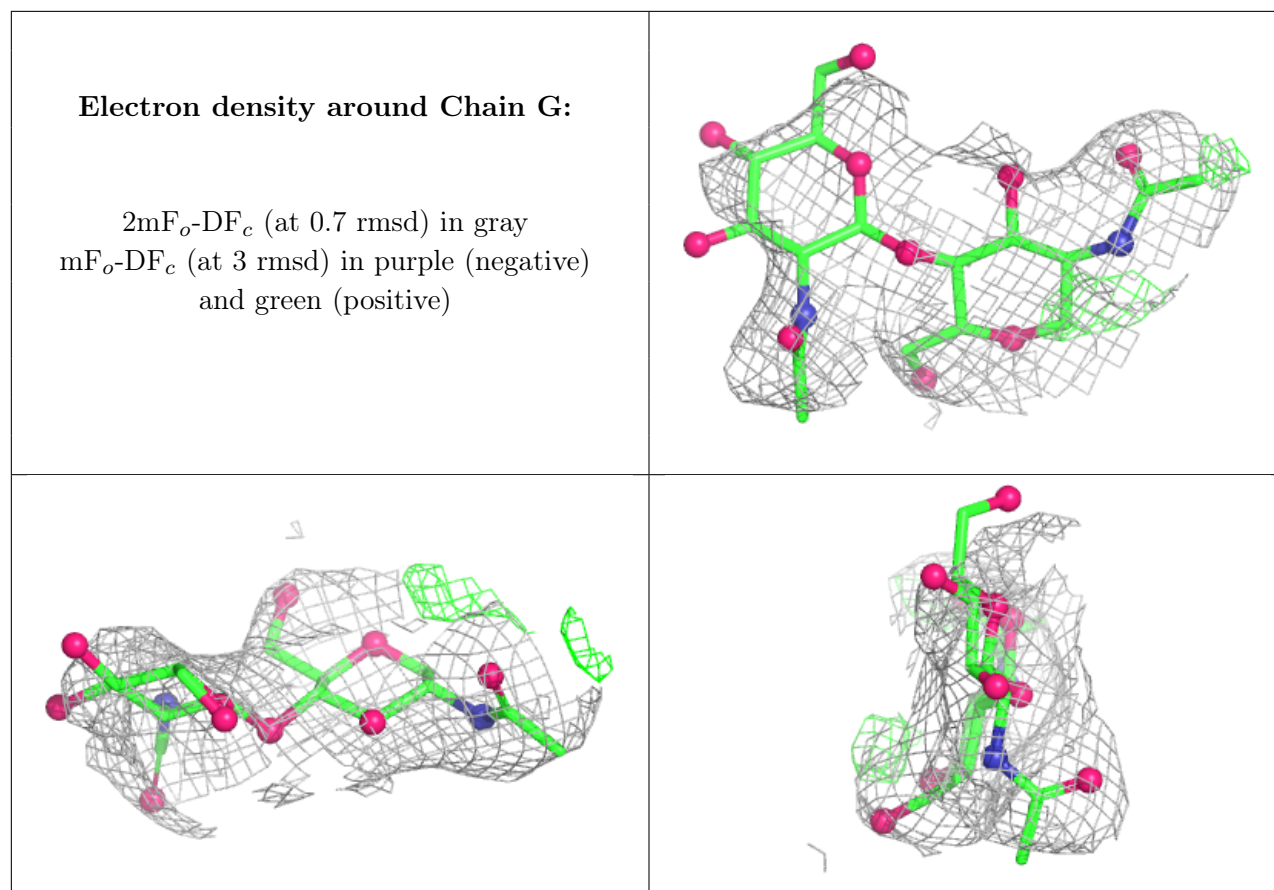
The following is a graphical depiction of the model fit to experimental electron density for oligosaccharide. Each fit is shown from different orientation to approximate a three-dimensional view.



Electron density around Chain F:

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





6.4 Ligands [i](#)

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
6	NAG	D	702	14/15	0.39	0.15	132,144,155,158	0
6	NAG	C	712	14/15	0.43	0.23	115,135,162,176	0
6	NAG	C	711	14/15	0.45	0.16	104,126,132,133	0
6	NAG	D	706	14/15	0.51	0.15	101,122,127,130	0
6	NAG	D	701	14/15	0.63	0.16	95,105,126,128	0
6	NAG	C	713	14/15	0.67	0.15	93,102,114,116	0
6	NAG	D	703	14/15	0.70	0.14	83,109,115,120	0

6.5 Other polymers [i](#)

There are no such residues in this entry.