



wwPDB X-ray Structure Validation Summary Report ⓘ

Aug 26, 2021 – 12:05 PM EDT

PDB ID : 4XQO
Title : Crystal structure of hemagglutinin from Jiangxi-Donghu (2013) H10N8 influenza virus in complex with 6'-SLN
Authors : Tzarum, N.; Zhang, H.; Zhu, X.; Wilson, I.A.
Deposited on : 2015-01-19
Resolution : 2.85 Å(reported)

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

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

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

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

MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.23.1
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)	:	2.23.1

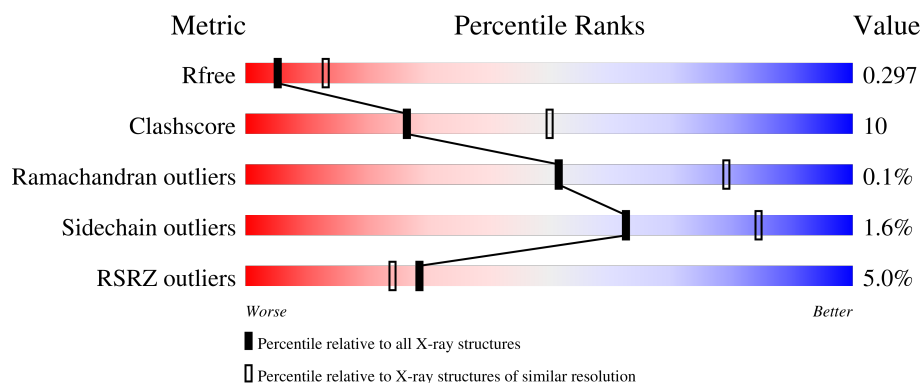
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.85 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	3168 (2.90-2.82)
Clashscore	141614	3438 (2.90-2.82)
Ramachandran outliers	138981	3348 (2.90-2.82)
Sidechain outliers	138945	3351 (2.90-2.82)
RSRZ outliers	127900	3103 (2.90-2.82)

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	326	<div> <div>3%</div> <div>79%</div> <div>17%</div> <div>..</div> </div>
1	C	326	<div> <div>3%</div> <div>74%</div> <div>22%</div> <div>..</div> </div>
1	E	326	<div> <div>3%</div> <div>74%</div> <div>23%</div> <div>.</div> </div>
2	B	181	<div> <div>7%</div> <div>64%</div> <div>29%</div> <div>• 7%</div> </div>
2	D	181	<div> <div>8%</div> <div>67%</div> <div>24%</div> <div>• 7%</div> </div>

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Mol	Chain	Length	Quality of chain
2	F	181	
3	G	2	
3	J	2	
4	H	3	
5	I	2	

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	NAG	H	1	-	-	-	X

2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 11538 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 Hemagglutinin HA1 chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	316	Total	C	N	O	S	0	0	0
			2424	1499	447	461	17			
1	C	315	Total	C	N	O	S	0	0	0
			2411	1491	445	458	17			
1	E	318	Total	C	N	O	S	0	0	0
			2437	1506	449	465	17			

There are 9 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	8	ALA	-	expression tag	UNP A0A059T4A1
A	9	ASP	-	expression tag	UNP A0A059T4A1
A	10	PRO	-	expression tag	UNP A0A059T4A1
C	8	ALA	-	expression tag	UNP A0A059T4A1
C	9	ASP	-	expression tag	UNP A0A059T4A1
C	10	PRO	-	expression tag	UNP A0A059T4A1
E	8	ALA	-	expression tag	UNP A0A059T4A1
E	9	ASP	-	expression tag	UNP A0A059T4A1
E	10	PRO	-	expression tag	UNP A0A059T4A1

- Molecule 2 is a protein called Hemagglutinin HA2 chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	169	Total	C	N	O	S	0	0	0
			1364	843	236	277	8			
2	D	168	Total	C	N	O	S	0	0	0
			1362	842	235	277	8			
2	F	166	Total	C	N	O	S	0	0	0
			1351	837	233	273	8			

There are 21 discrepancies between the modelled and reference sequences:

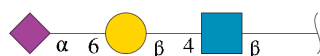
Chain	Residue	Modelled	Actual	Comment	Reference
B	175	SER	-	expression tag	UNP A0A059T4A1
B	176	GLY	-	expression tag	UNP A0A059T4A1
B	177	ARG	-	expression tag	UNP A0A059T4A1
B	178	LEU	-	expression tag	UNP A0A059T4A1
B	179	VAL	-	expression tag	UNP A0A059T4A1
B	180	PRO	-	expression tag	UNP A0A059T4A1
B	181	ARG	-	expression tag	UNP A0A059T4A1
D	175	SER	-	expression tag	UNP A0A059T4A1
D	176	GLY	-	expression tag	UNP A0A059T4A1
D	177	ARG	-	expression tag	UNP A0A059T4A1
D	178	LEU	-	expression tag	UNP A0A059T4A1
D	179	VAL	-	expression tag	UNP A0A059T4A1
D	180	PRO	-	expression tag	UNP A0A059T4A1
D	181	ARG	-	expression tag	UNP A0A059T4A1
F	175	SER	-	expression tag	UNP A0A059T4A1
F	176	GLY	-	expression tag	UNP A0A059T4A1
F	177	ARG	-	expression tag	UNP A0A059T4A1
F	178	LEU	-	expression tag	UNP A0A059T4A1
F	179	VAL	-	expression tag	UNP A0A059T4A1
F	180	PRO	-	expression tag	UNP A0A059T4A1
F	181	ARG	-	expression tag	UNP A0A059T4A1

- Molecule 3 is an oligosaccharide called 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.



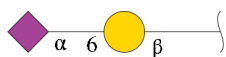
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
3	G	2	Total	C	N	O	0	0	0
			28	16	2	10			
3	J	2	Total	C	N	O	0	0	0
			28	16	2	10			

- Molecule 4 is an oligosaccharide called N-acetyl-alpha-neuraminic acid-(2-6)-beta-D-galactopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.



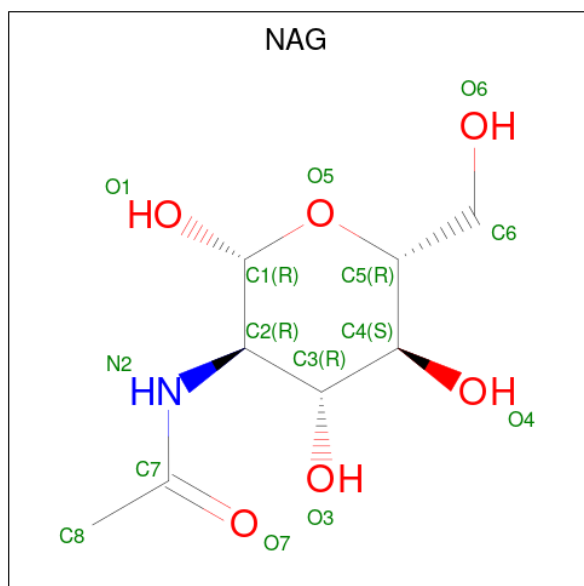
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
4	H	3	Total	C	N	O	0	0	0
			46	25	2	19			

- Molecule 5 is an oligosaccharide called N-acetyl-alpha-neuraminic acid-(2-6)-beta-D-galactopyranose.



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
5	I	2	Total	C	N	O	0	0	0
			32	17	1	14			

- Molecule 6 is 2-acetamido-2-deoxy-beta-D-glucopyranose (three-letter code: NAG) (formula: $C_8H_{15}NO_6$).



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

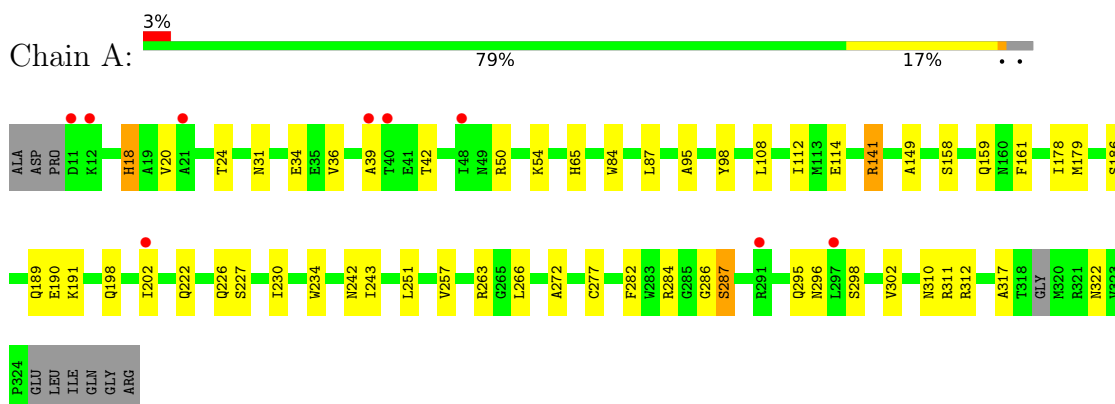
- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	12	Total 12	O 12	0	0
7	B	4	Total 4	O 4	0	0
7	C	6	Total 6	O 6	0	0
7	D	1	Total 1	O 1	0	0
7	E	4	Total 4	O 4	0	0

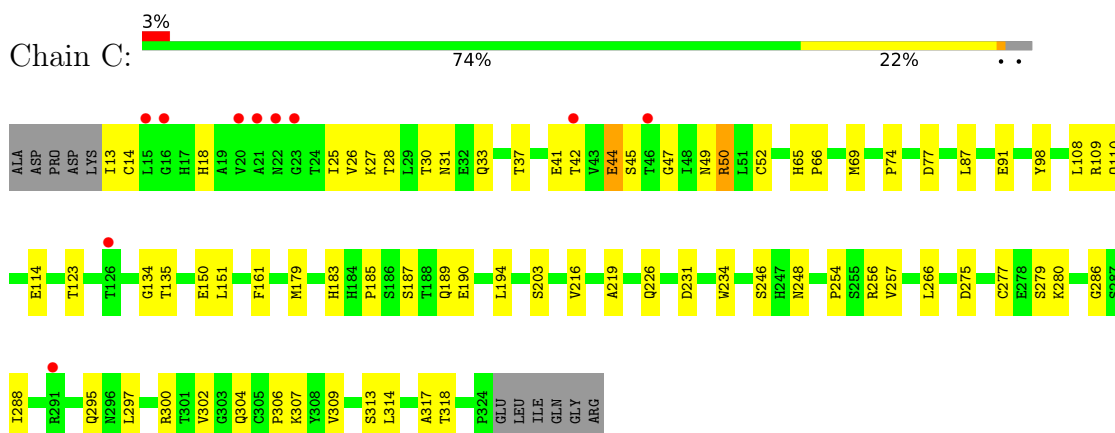
3 Residue-property plots [i](#)

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.

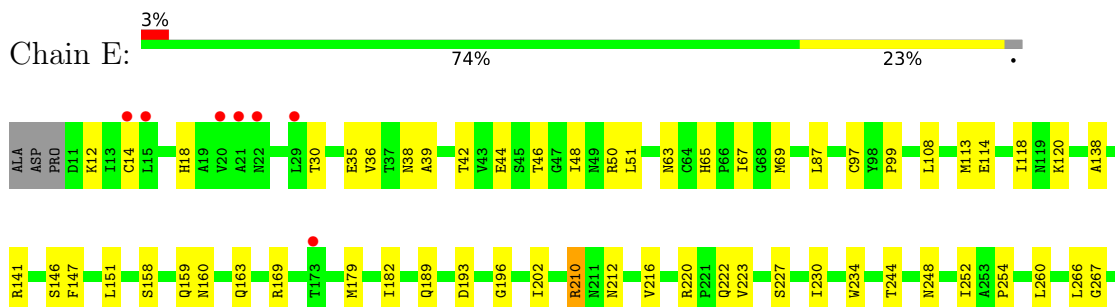
• Molecule 1: Hemagglutinin HA1 chain



• Molecule 1: Hemagglutinin HA1 chain

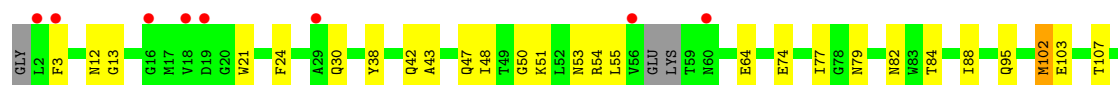


• Molecule 1: Hemagglutinin HA1 chain





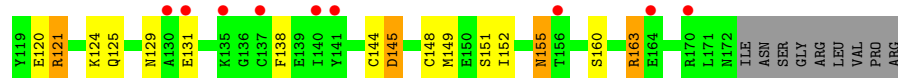
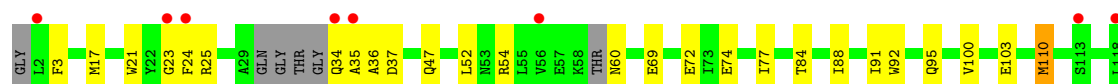
- Molecule 2: Hemagglutinin HA2 chain



- Molecule 2: Hemagglutinin HA2 chain



- Molecule 2: Hemagglutinin HA2 chain



- Molecule 3: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose



- Molecule 3: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose





- Molecule 4: N-acetyl-alpha-neuraminic acid-(2-6)-beta-D-galactopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain H:  67% 33%



- Molecule 5: N-acetyl-alpha-neuraminic acid-(2-6)-beta-D-galactopyranose

Chain I:  100%



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	64.07Å 245.42Å 71.30Å 90.00° 112.46° 90.00°	Depositor
Resolution (Å)	46.17 – 2.85 46.17 – 2.81	Depositor EDS
% Data completeness (in resolution range)	88.3 (46.17-2.85) 80.8 (46.17-2.81)	Depositor EDS
R_{merge}	0.14	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.43 (at 2.81Å)	Xtriage
Refinement program	REFMAC	Depositor
R, R_{free}	0.259 , 0.291 0.279 , 0.297	Depositor DCC
R_{free} test set	2125 reflections (5.01%)	wwPDB-VP
Wilson B-factor (Å ²)	56.9	Xtriage
Anisotropy	0.563	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.26 , 29.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.84	EDS
Total number of atoms	11538	wwPDB-VP
Average B, all atoms (Å ²)	68.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.75% 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, SIA, GAL

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	0.24	0/2472	0.48	0/3348
1	C	0.24	0/2460	0.46	0/3334
1	E	0.23	0/2486	0.45	0/3368
2	B	0.26	0/1388	0.49	0/1873
2	D	0.26	0/1385	0.47	0/1867
2	F	0.27	0/1374	0.48	0/1851
All	All	0.25	0/11565	0.47	0/15641

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
2	F	0	1

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	F	155	ASN	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	2424	0	2375	36	0
1	C	2411	0	2363	63	0
1	E	2437	0	2386	52	0
2	B	1364	0	1259	43	0
2	D	1362	0	1254	42	0
2	F	1351	0	1249	32	0
3	G	28	0	25	1	0
3	J	28	0	25	1	0
4	H	46	0	40	2	0
5	I	32	0	27	4	0
6	A	14	0	13	0	0
6	C	14	0	13	0	0
7	A	12	0	0	3	0
7	B	4	0	0	1	0
7	C	6	0	0	0	0
7	D	1	0	0	0	0
7	E	4	0	0	1	0
All	All	11538	0	11029	232	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 10.

The worst 5 of 232 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:C:44:GLU:OE1	1:C:45:SER:N	2.11	0.80
2:B:125:GLN:HE21	2:B:157:TYR:HB3	1.48	0.78
1:A:226:GLN:NE2	4:H:3:SIA:O1A	2.17	0.77
1:E:63:ASN:OD1	7:E:701:HOH:O	2.03	0.76
1:C:30:THR:HG22	2:D:105:GLN:HE22	1.51	0.74

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	312/326 (96%)	302 (97%)	10 (3%)	0	100	100
1	C	313/326 (96%)	295 (94%)	18 (6%)	0	100	100
1	E	316/326 (97%)	305 (96%)	10 (3%)	1 (0%)	41	68
2	B	165/181 (91%)	161 (98%)	4 (2%)	0	100	100
2	D	162/181 (90%)	149 (92%)	12 (7%)	1 (1%)	25	53
2	F	160/181 (88%)	157 (98%)	3 (2%)	0	100	100
All	All	1428/1521 (94%)	1369 (96%)	57 (4%)	2 (0%)	51	79

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	D	144	CYS
1	E	324	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	268/275 (98%)	264 (98%)	4 (2%)	65	86
1	C	266/275 (97%)	264 (99%)	2 (1%)	81	93
1	E	269/275 (98%)	266 (99%)	3 (1%)	73	90
2	B	144/154 (94%)	143 (99%)	1 (1%)	84	94
2	D	144/154 (94%)	141 (98%)	3 (2%)	53	79
2	F	143/154 (93%)	136 (95%)	7 (5%)	25	54
All	All	1234/1287 (96%)	1214 (98%)	20 (2%)	62	84

5 of 20 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	F	74	GLU

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Mol	Chain	Res	Type
2	F	144	CYS
2	F	163	ARG
2	F	145	ASP
1	C	50	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 36 such sidechains are listed below:

Mol	Chain	Res	Type
2	F	60	ASN
2	F	172	ASN
2	F	117	ASN
2	F	155	ASN
1	C	226	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 ⓘ

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	$\# Z > 2$
3	NAG	G	1	3,1	14,14,15	0.76	1 (7%)	17,19,21	0.89	0
3	NAG	G	2	3	14,14,15	0.45	0	17,19,21	0.92	1 (5%)
4	NAG	H	1	4	15,15,15	0.47	0	21,21,21	1.48	4 (19%)
4	GAL	H	2	4	11,11,12	0.45	0	15,15,17	0.99	1 (6%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
4	SIA	H	3	4	17,20,21	0.36	0	21,28,31	1.16	3 (14%)
5	GAL	I	1	5	12,12,12	1.10	1 (8%)	17,17,17	0.89	1 (5%)
5	SIA	I	2	5	17,20,21	0.31	0	21,28,31	1.17	2 (9%)
3	NAG	J	1	3,1	14,14,15	0.75	1 (7%)	17,19,21	0.95	1 (5%)
3	NAG	J	2	3	14,14,15	0.33	0	17,19,21	0.78	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	G	1	3,1	-	0/6/23/26	0/1/1/1
3	NAG	G	2	3	-	4/6/23/26	0/1/1/1
4	NAG	H	1	4	-	1/6/26/26	0/1/1/1
4	GAL	H	2	4	-	0/2/19/22	0/1/1/1
4	SIA	H	3	4	-	0/14/34/38	0/1/1/1
5	GAL	I	1	5	-	0/2/22/22	0/1/1/1
5	SIA	I	2	5	-	0/14/34/38	0/1/1/1
3	NAG	J	1	3,1	-	2/6/23/26	0/1/1/1
3	NAG	J	2	3	-	0/6/23/26	0/1/1/1

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	I	1	GAL	O1-C1	-3.68	1.27	1.39
3	J	1	NAG	O5-C1	-2.48	1.39	1.43
3	G	1	NAG	O5-C1	-2.43	1.39	1.43

The worst 5 of 13 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	H	1	NAG	C1-C2-N2	-3.39	106.80	110.73
4	H	1	NAG	C4-C3-C2	3.35	115.25	110.34
5	I	2	SIA	C4-C5-N5	-3.15	104.13	110.38
4	H	1	NAG	C3-C4-C5	2.76	115.16	110.24
4	H	3	SIA	C6-O6-C2	2.73	117.17	111.34

There are no chirality outliers.

5 of 7 torsion outliers are listed below:

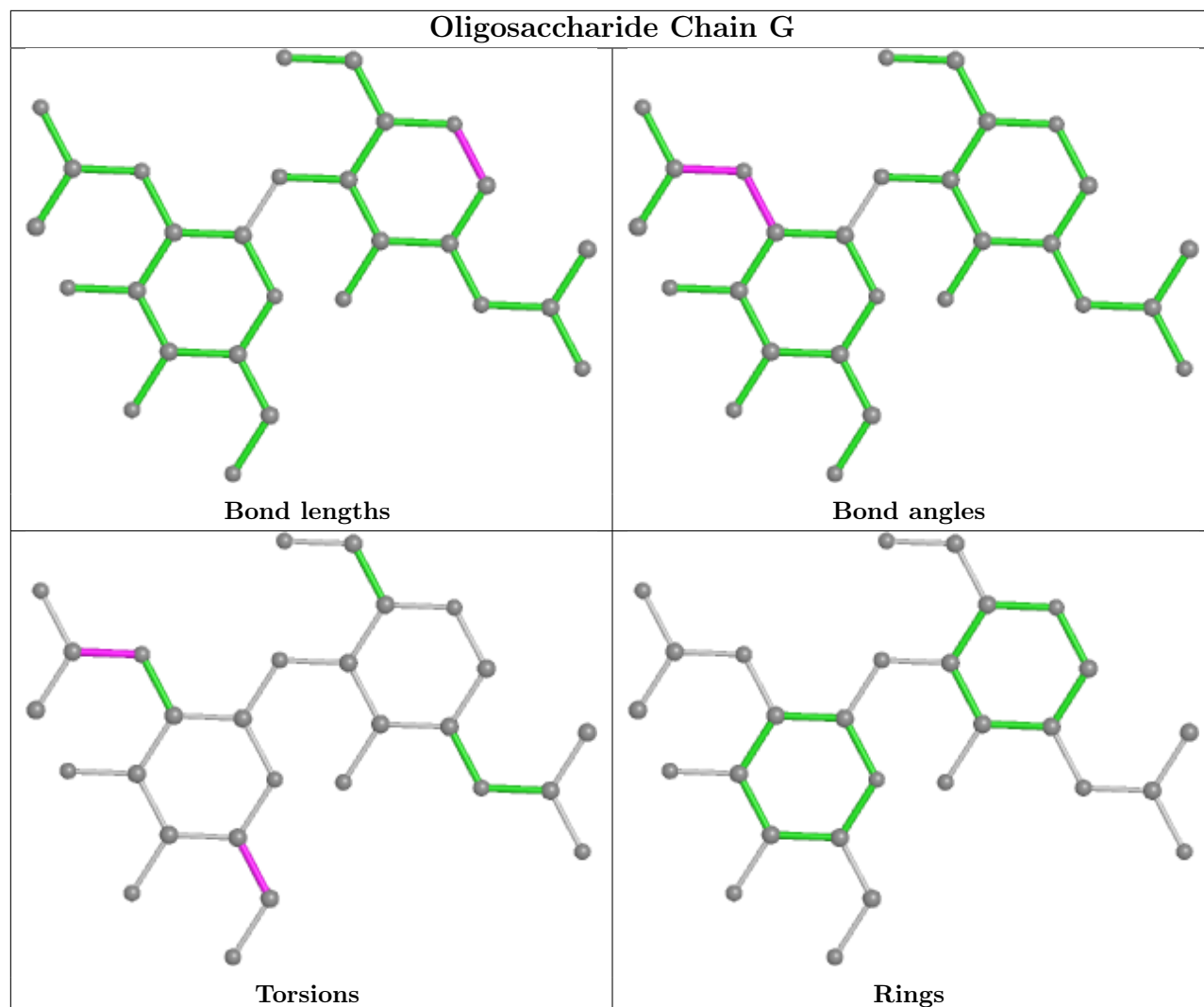
Mol	Chain	Res	Type	Atoms
3	G	2	NAG	O5-C5-C6-O6
3	J	1	NAG	O5-C5-C6-O6
3	G	2	NAG	C4-C5-C6-O6
3	J	1	NAG	C4-C5-C6-O6
3	G	2	NAG	C8-C7-N2-C2

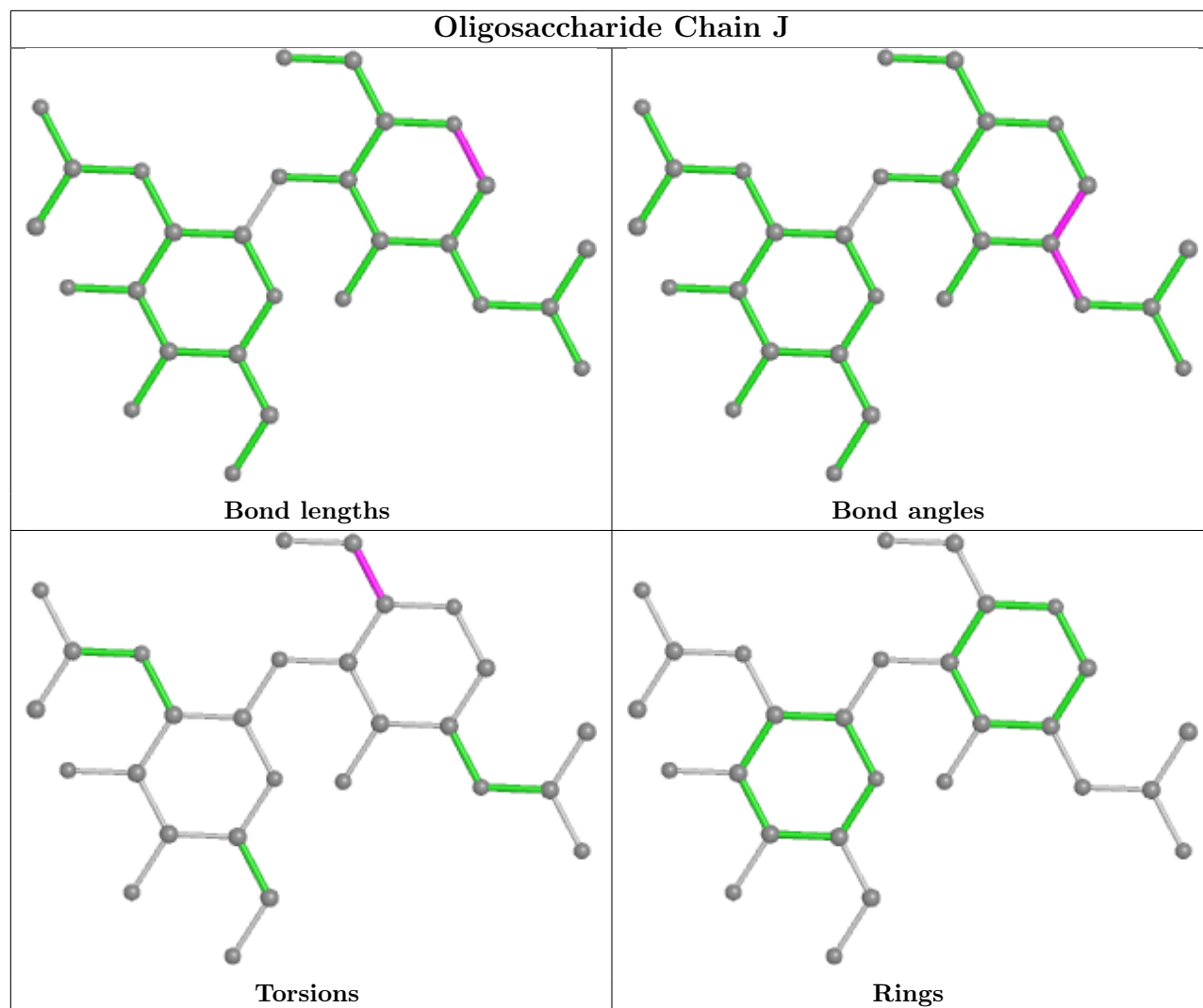
There are no ring outliers.

6 monomers are involved in 8 short contacts:

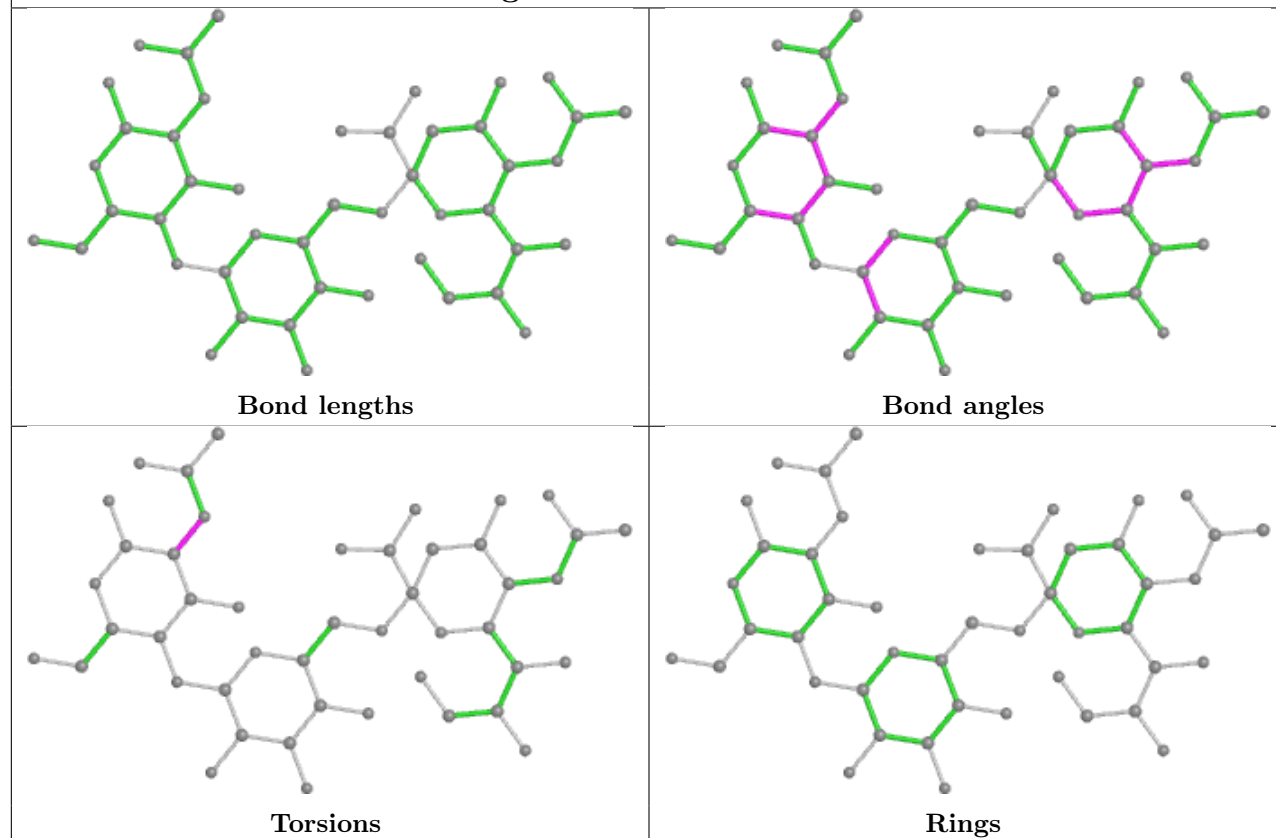
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	G	2	NAG	1	0
5	I	2	SIA	3	0
5	I	1	GAL	1	0
3	G	1	NAG	1	0
3	J	1	NAG	1	0
4	H	3	SIA	2	0

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

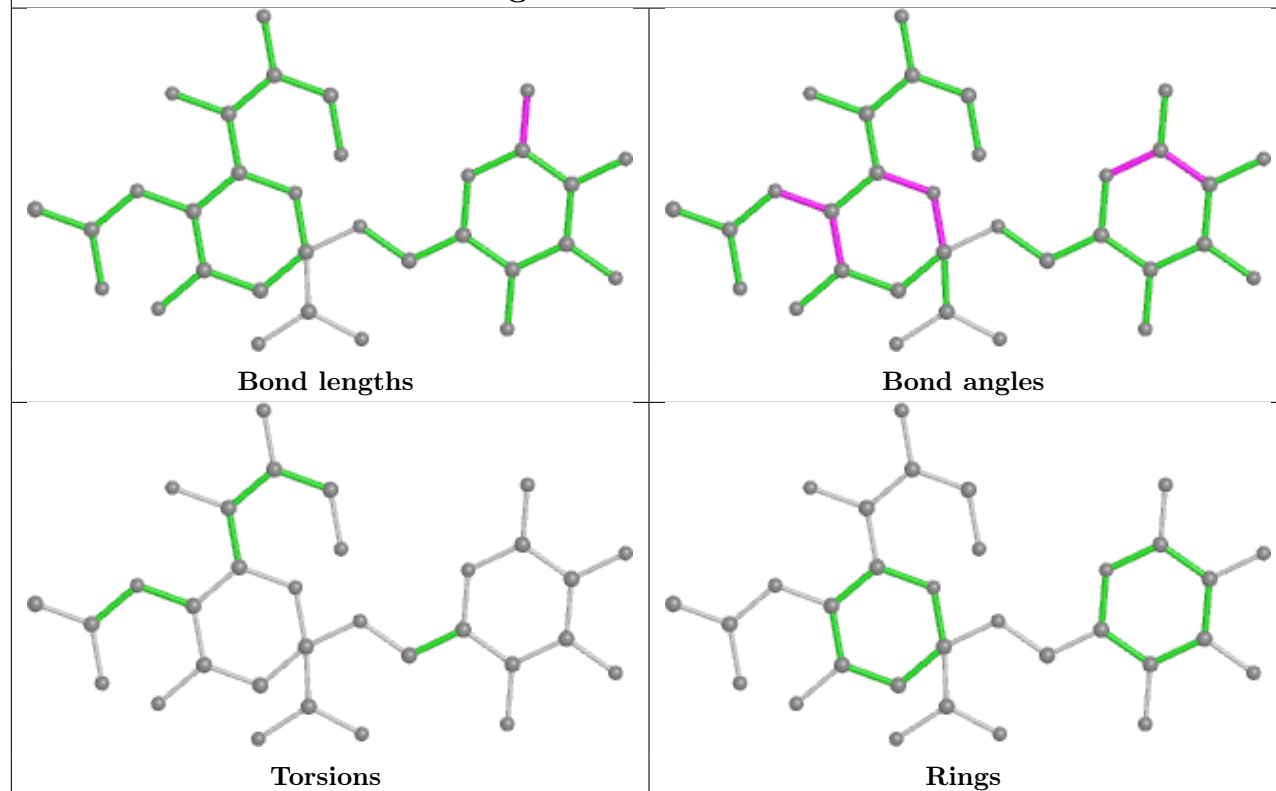




Oligosaccharide Chain H



Oligosaccharide Chain I



5.6 Ligand geometry

2 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	A	401	1	14,14,15	0.65	1 (7%)	17,19,21	0.57	0
6	NAG	C	501	1	14,14,15	0.79	1 (7%)	17,19,21	0.79	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	A	401	1	-	0/6/23/26	0/1/1/1
6	NAG	C	501	1	-	2/6/23/26	0/1/1/1

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	C	501	NAG	O5-C1	-2.45	1.39	1.43
6	A	401	NAG	O5-C1	-2.35	1.40	1.43

There are no bond angle outliers.

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
6	C	501	NAG	C4-C5-C6-O6
6	C	501	NAG	O5-C5-C6-O6

There are no ring outliers.

No monomer is involved in short contacts.

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	316/326 (96%)	0.12	9 (2%) 53 48	39, 53, 82, 99	0
1	C	315/326 (96%)	0.24	10 (3%) 47 42	40, 59, 87, 101	0
1	E	318/326 (97%)	0.20	10 (3%) 49 44	37, 60, 98, 113	0
2	B	169/181 (93%)	0.60	13 (7%) 13 10	45, 81, 100, 108	0
2	D	168/181 (92%)	0.68	14 (8%) 11 8	53, 83, 103, 112	0
2	F	166/181 (91%)	0.86	17 (10%) 6 4	42, 88, 108, 117	0
All	All	1452/1521 (95%)	0.37	73 (5%) 28 24	37, 65, 100, 117	0

The worst 5 of 73 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	F	23	GLY	7.0
1	E	14	CYS	4.8
2	D	8	GLY	4.4
2	F	170	ARG	4.4
1	C	16	GLY	4.2

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

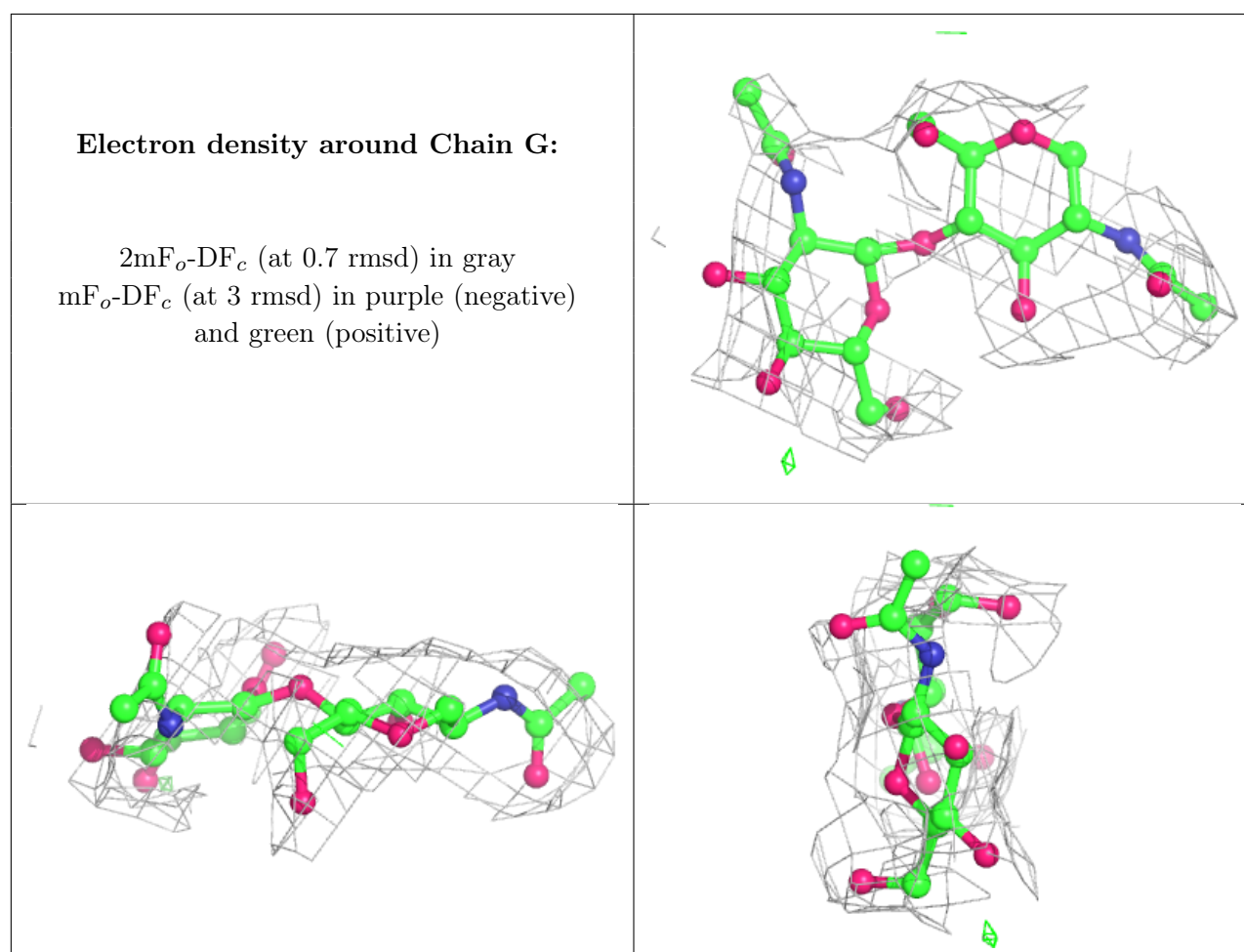
There are no non-standard protein/DNA/RNA residues in this entry.

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.

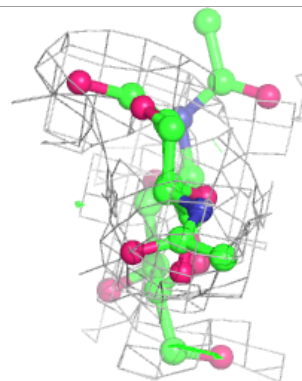
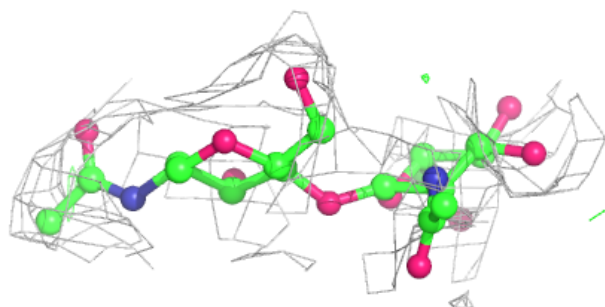
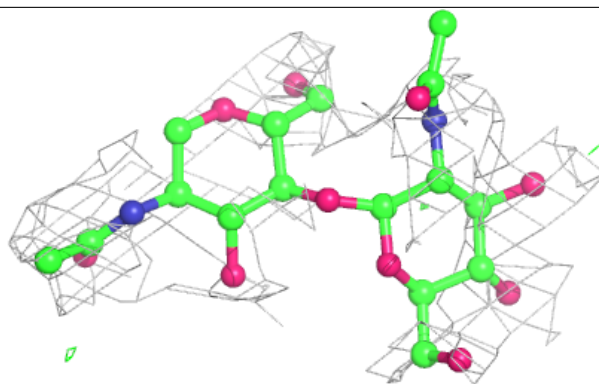
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
3	NAG	J	2	14/15	0.56	0.38	123,128,143,144	0
3	NAG	G	2	14/15	0.65	0.31	104,113,122,125	0
4	NAG	H	1	15/15	0.72	0.43	88,104,109,117	0
4	GAL	H	2	11/12	0.77	0.25	71,82,94,97	0
4	SIA	H	3	20/21	0.81	0.23	54,72,82,83	0
3	NAG	G	1	14/15	0.82	0.18	64,83,92,101	0
5	GAL	I	1	12/12	0.82	0.17	68,92,104,105	0
3	NAG	J	1	14/15	0.87	0.15	74,86,102,105	0
5	SIA	I	2	20/21	0.91	0.18	50,63,79,81	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.



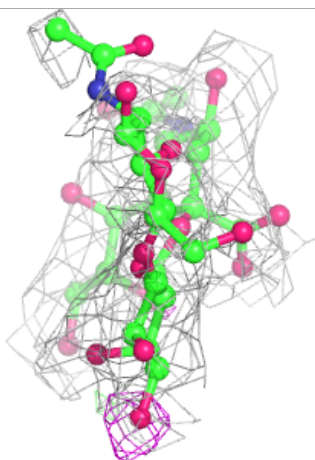
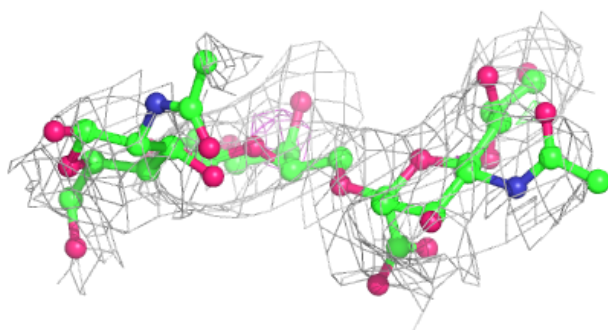
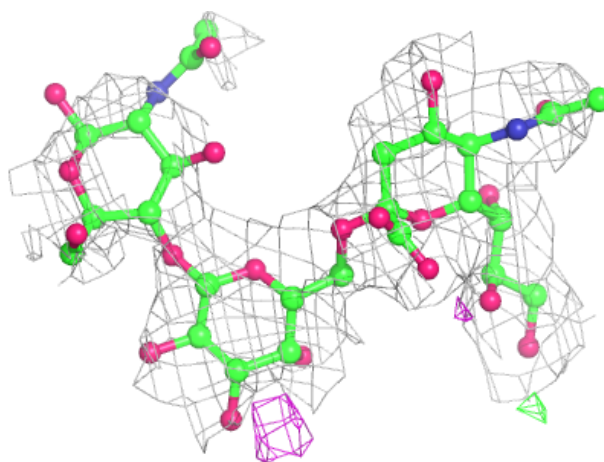
Electron density around Chain J:

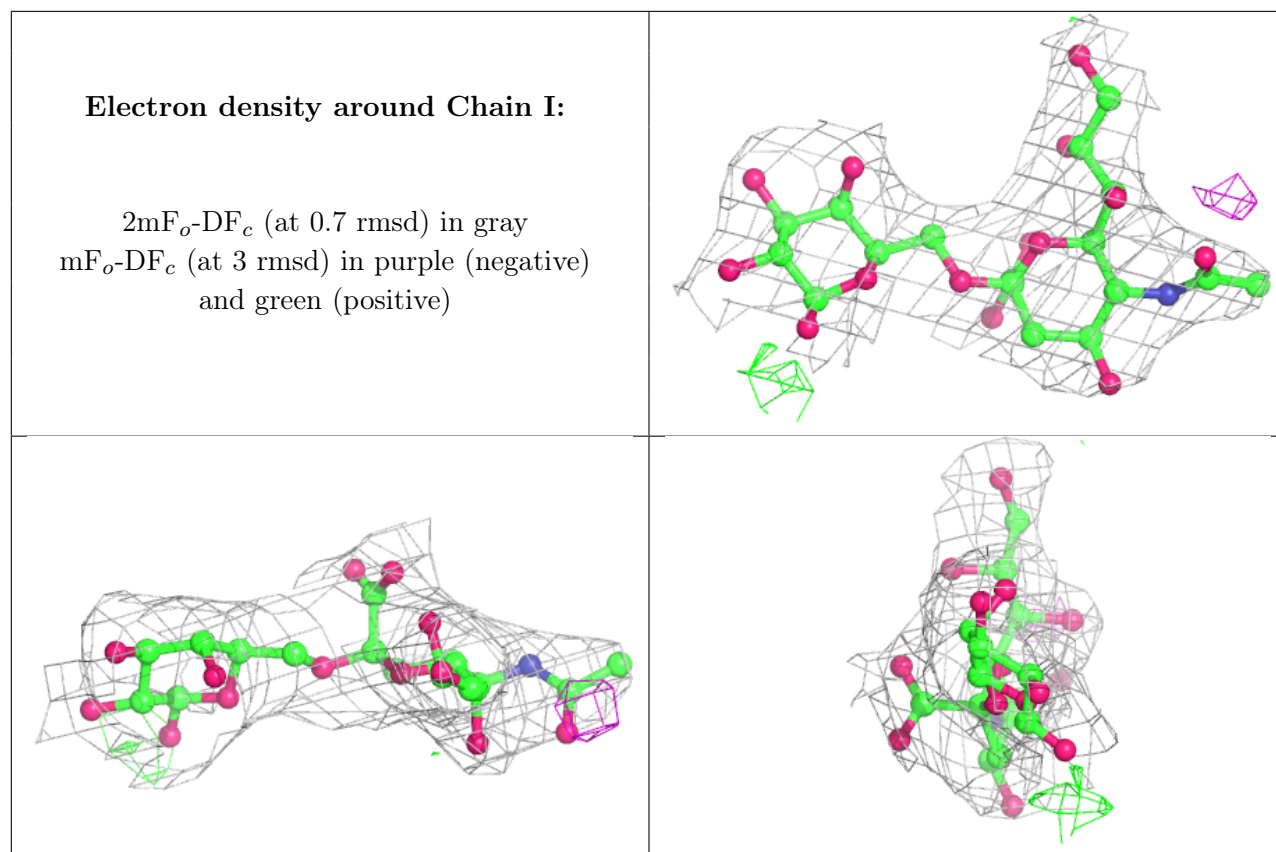
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around Chain H:

$2mF_o-DF_c$ (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(\AA^2)	Q<0.9
6	NAG	A	401	14/15	0.66	0.28	86,98,107,112	0
6	NAG	C	501	14/15	0.81	0.18	61,80,94,96	0

6.5 Other polymers [i](#)

There are no such residues in this entry.