wwPDB X-ray Structure Validation Summary Report (i)

Jun 22, 2024 – 10:10 AM EDT

PDB ID	:	4QY1
Title	:	Structure of H10 from human-infecting H10N8 in complex with avian receptor
Authors	:	Wang, M.; Zhang, W.; Qi, J.; Wang, F.; Zhou, J.; Bi, Y.; Wu, Y.; Sun, H.; Liu, J.; Huang, C.; Li, X.; Yan, J.; Shu, Y.; Shi, Y.; Gao, G.F.
Deposited on	:	2014-07-23
Resolution	:	2.59 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 (i) 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 (1)) were used in the production of this report:

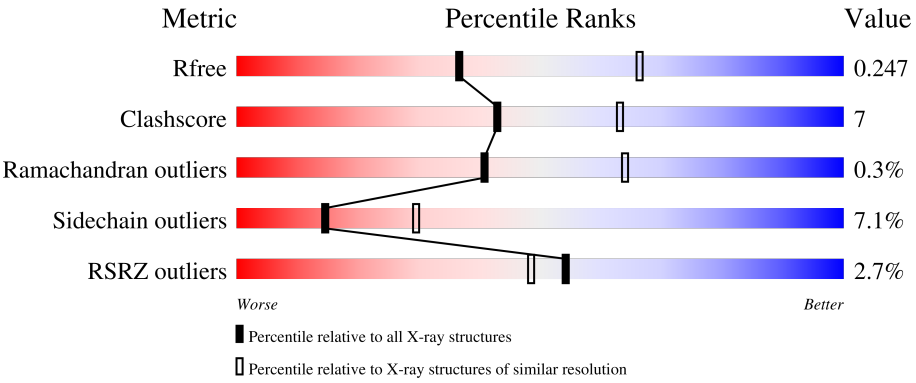
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.37.1
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.37.1

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.59 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	3163 (2.60-2.60)
Clashscore	141614	3518 (2.60-2.60)
Ramachandran outliers	138981	3455 (2.60-2.60)
Sidechain outliers	138945	3455 (2.60-2.60)
RSRZ outliers	127900	3104 (2.60-2.60)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for >=3, 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions <=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	318	78% 20%
1	C	318	% 81% 17%
1	E	318	79% 19%
1	G	318	% 81% 18%
1	I	318	2% 82% 15%

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Mol	Chain	Length	Quality of chain
1	K	318	2% 81% 18%
1	M	318	10% 83% 15%
1	O	318	8% 78% 21%
1	Q	318	% 76% 22%
1	S	318	6% 82% 16%
1	U	318	2% 81% 16%
1	W	318	3% 79% 19%
2	B	174	2% 83% 14%
2	D	174	2% 82% 16%
2	F	174	2% 82% 16%
2	H	174	% 82% 16%
2	J	174	3% 86% 13%
2	L	174	3% 84% 14%
2	N	174	2% 84% 14%
2	P	174	3% 82% 16%
2	R	174	5% 82% 17%
2	T	174	2% 82% 16%
2	V	174	2% 80% 17%
2	X	174	2% 82% 17%
3	Y	2	50% 50%
4	Z	3	67% 33%
4	a	3	100%
4	b	3	67% 33%
4	c	3	33% 67%

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 crite-

ria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
4	NAG	Z	1	-	-	-	X
4	NAG	a	1	-	-	-	X
4	NAG	b	1	-	-	-	X
4	NAG	c	1	-	-	-	X

2 Entry composition (i)

There are 6 unique types of molecules in this entry. The entry contains 48138 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.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	318	Total	C	N	O	S	0	0	0
			2437	1506	449	465	17			
1	C	318	Total	C	N	O	S	0	0	0
			2437	1506	449	465	17			
1	E	318	Total	C	N	O	S	0	0	0
			2437	1506	449	465	17			
1	G	318	Total	C	N	O	S	0	0	0
			2437	1506	449	465	17			
1	I	318	Total	C	N	O	S	0	0	0
			2437	1506	449	465	17			
1	K	318	Total	C	N	O	S	0	0	0
			2437	1506	449	465	17			
1	M	318	Total	C	N	O	S	0	0	0
			2437	1506	449	465	17			
1	O	318	Total	C	N	O	S	0	0	0
			2437	1506	449	465	17			
1	Q	318	Total	C	N	O	S	0	0	0
			2437	1506	449	465	17			
1	S	318	Total	C	N	O	S	0	0	0
			2436	1506	449	464	17			
1	U	318	Total	C	N	O	S	0	0	0
			2437	1506	449	465	17			
1	W	318	Total	C	N	O	S	0	0	0
			2437	1506	449	465	17			

- Molecule 2 is a protein called hemagglutinin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	174	Total	C	N	O	S	0	0	0
			1402	866	243	285	8			
2	D	174	Total	C	N	O	S	0	0	0
			1402	866	243	285	8			

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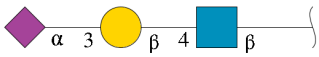
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	F	174	Total	C	N	O	S	0	0	0
			1402	866	243	285	8			
2	H	174	Total	C	N	O	S	0	0	0
			1402	866	243	285	8			
2	J	174	Total	C	N	O	S	0	0	0
			1402	866	243	285	8			
2	L	174	Total	C	N	O	S	0	0	0
			1402	866	243	285	8			
2	N	174	Total	C	N	O	S	0	0	0
			1402	866	243	285	8			
2	P	174	Total	C	N	O	S	0	0	0
			1402	866	243	285	8			
2	R	174	Total	C	N	O	S	0	0	0
			1402	866	243	285	8			
2	T	174	Total	C	N	O	S	0	0	0
			1402	866	243	285	8			
2	V	174	Total	C	N	O	S	0	0	0
			1402	866	243	285	8			
2	X	174	Total	C	N	O	S	0	0	0
			1402	866	243	285	8			

- 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	Y	2	Total	C	N	O	0	0	0
			28	16	2	10			

- Molecule 4 is an oligosaccharide called N-acetyl-alpha-neuraminic acid-(2-3)-beta-D-galactopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
4	Z	3	Total	C	N	O	0	0	0
			45	25	2	18			

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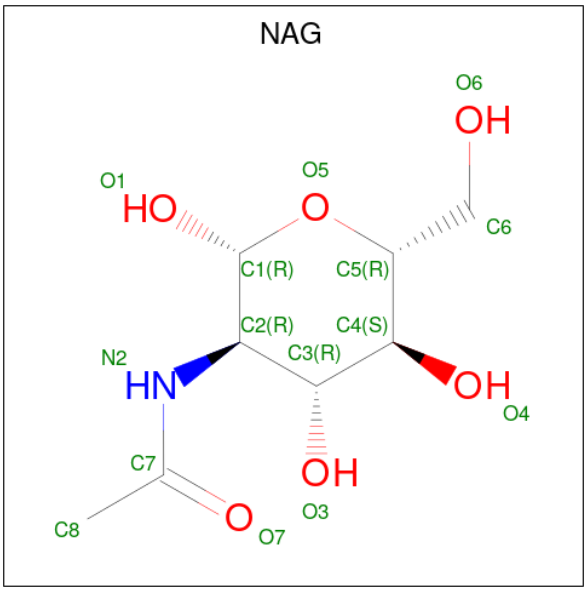
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf	Trace
4	a	3	Total C N O 45 25 2 18	0	0	0
4	b	3	Total C N O 45 25 2 18	0	0	0
4	c	3	Total C N O 45 25 2 18	0	0	0

- Molecule 5 is 2-acetamido-2-deoxy-beta-D-glucopyranose (three-letter code: NAG) (formula: $C_8H_{15}NO_6$).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	1	Total C N O 14 8 1 5	0	0
5	B	1	Total C N O 14 8 1 5	0	0
5	C	1	Total C N O 14 8 1 5	0	0
5	D	1	Total C N O 14 8 1 5	0	0
5	E	1	Total C N O 14 8 1 5	0	0
5	E	1	Total C N O 14 8 1 5	0	0
5	F	1	Total C N O 14 8 1 5	0	0
5	G	1	Total C N O 14 8 1 5	0	0

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Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	H	1	Total C N O 14 8 1 5	0	0
5	I	1	Total C N O 14 8 1 5	0	0
5	J	1	Total C N O 14 8 1 5	0	0
5	K	1	Total C N O 14 8 1 5	0	0
5	L	1	Total C N O 14 8 1 5	0	0
5	M	1	Total C N O 14 8 1 5	0	0
5	N	1	Total C N O 14 8 1 5	0	0
5	O	1	Total C N O 14 8 1 5	0	0
5	P	1	Total C N O 14 8 1 5	0	0
5	Q	1	Total C N O 14 8 1 5	0	0
5	R	1	Total C N O 14 8 1 5	0	0
5	S	1	Total C N O 14 8 1 5	0	0
5	T	1	Total C N O 14 8 1 5	0	0
5	U	1	Total C N O 14 8 1 5	0	0
5	V	1	Total C N O 14 8 1 5	0	0
5	W	1	Total C N O 14 8 1 5	0	0
5	X	1	Total C N O 14 8 1 5	0	0

- Molecule 6 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	A	106	Total O 106 106	0	0
6	B	70	Total O 70 70	0	0

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Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	C	71	Total O 71 71	0	0
6	D	87	Total O 87 87	0	0
6	E	93	Total O 93 93	0	0
6	F	64	Total O 64 64	0	0
6	G	80	Total O 80 80	0	0
6	H	53	Total O 53 53	0	0
6	I	102	Total O 102 102	0	0
6	J	43	Total O 43 43	0	0
6	K	93	Total O 93 93	0	0
6	L	53	Total O 53 53	0	0
6	M	27	Total O 27 27	0	0
6	N	46	Total O 46 46	0	0
6	O	40	Total O 40 40	0	0
6	P	48	Total O 48 48	0	0
6	Q	84	Total O 84 84	0	0
6	R	42	Total O 42 42	0	0
6	S	64	Total O 64 64	0	0
6	T	49	Total O 49 49	0	0
6	U	76	Total O 76 76	0	0
6	V	44	Total O 44 44	0	0
6	W	50	Total O 50 50	0	0

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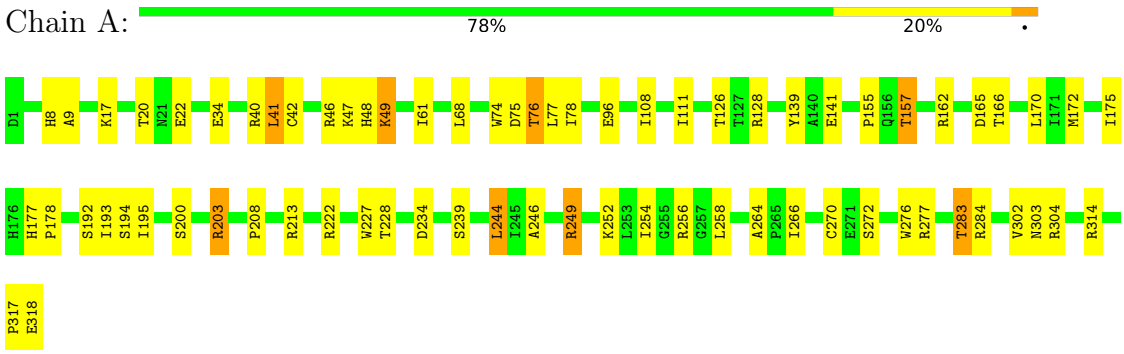
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	X	28	Total 28	O 28	0	0

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.

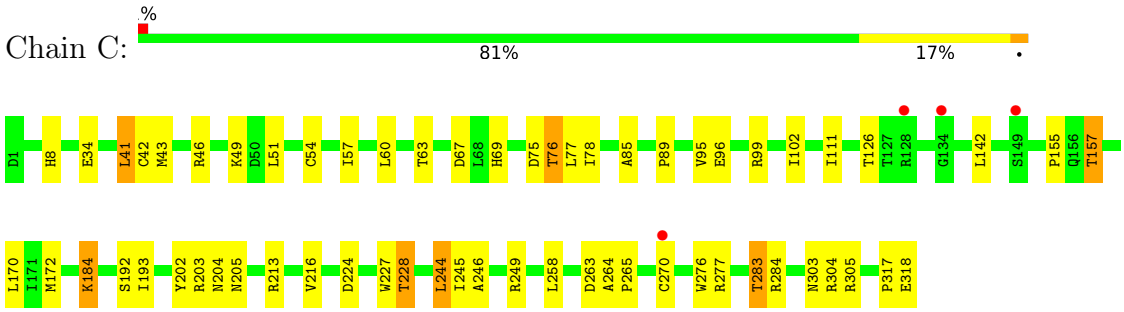
• Molecule 1: hemagglutinin

Chain A:



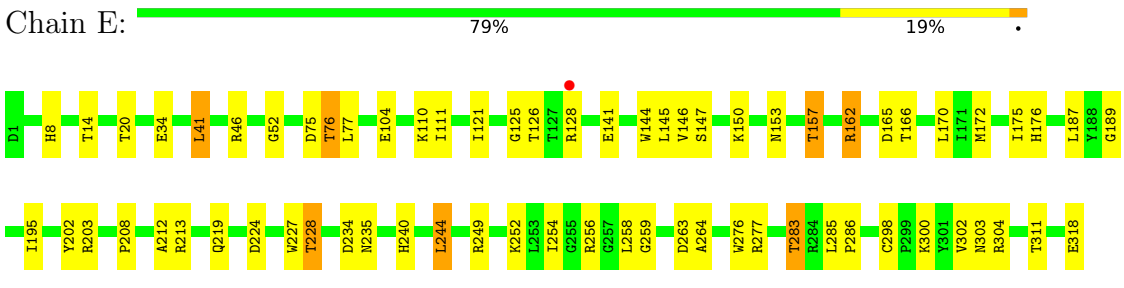
• Molecule 1: hemagglutinin

Chain C:



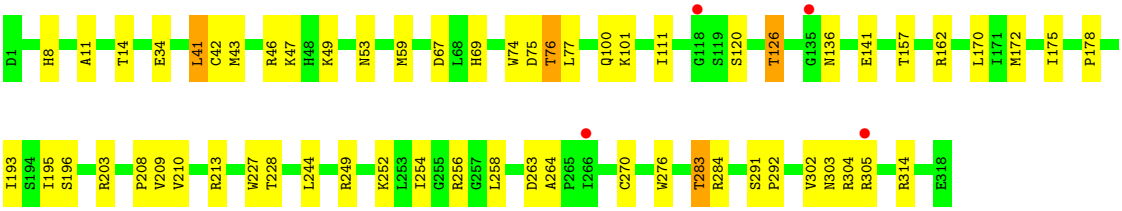
• Molecule 1: hemagglutinin

Chain E:

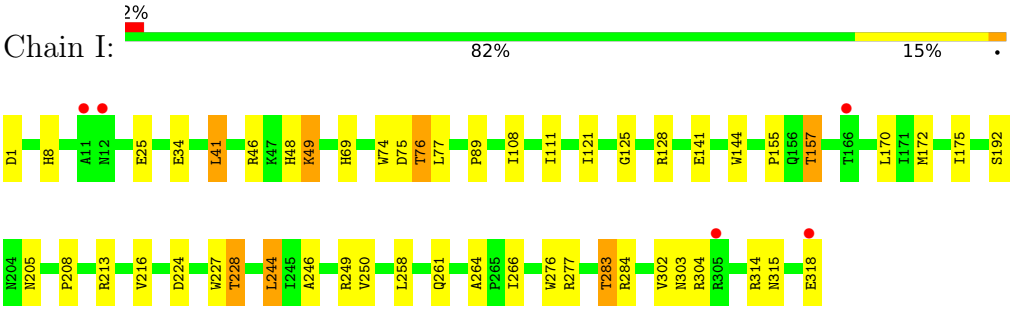


• Molecule 1: hemagglutinin

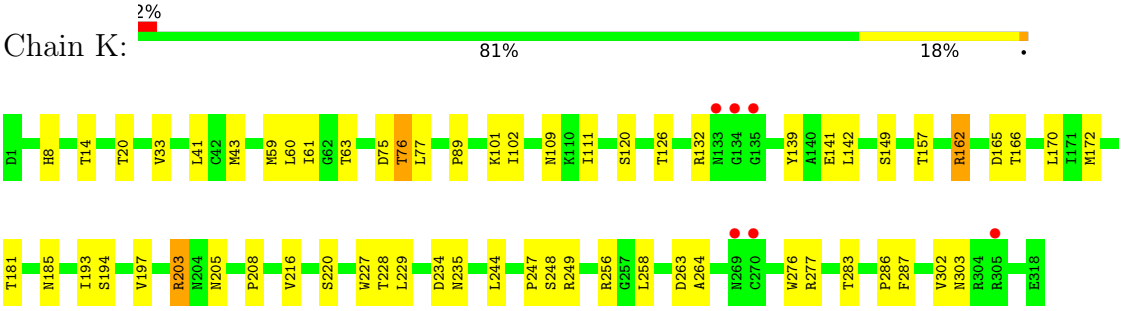
Chain G:



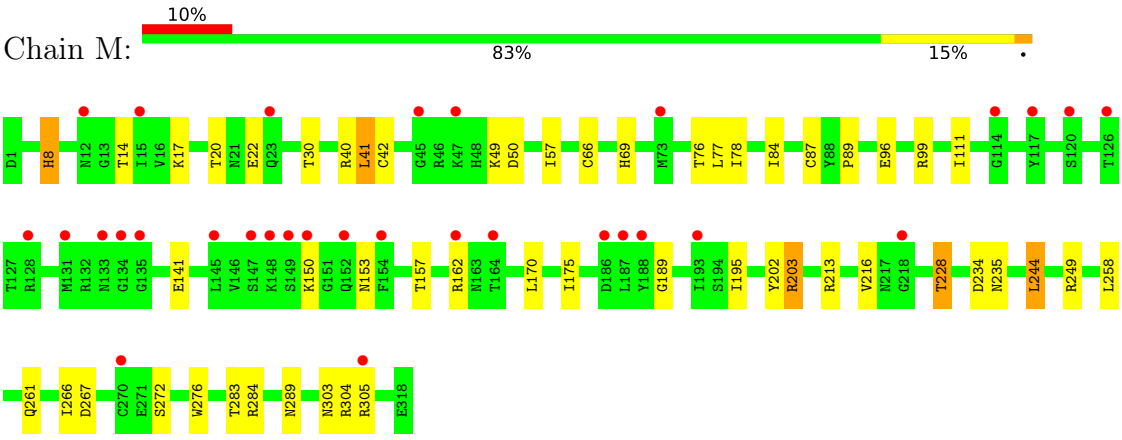
• Molecule 1: hemagglutinin



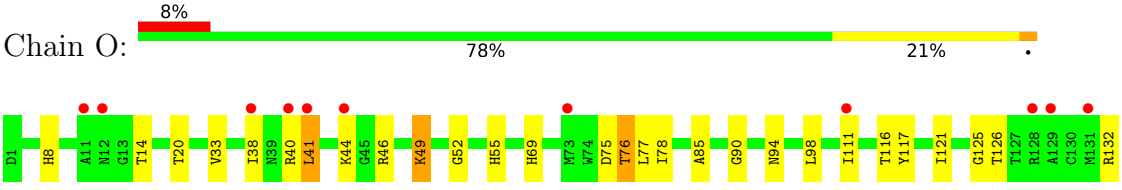
• Molecule 1: hemagglutinin

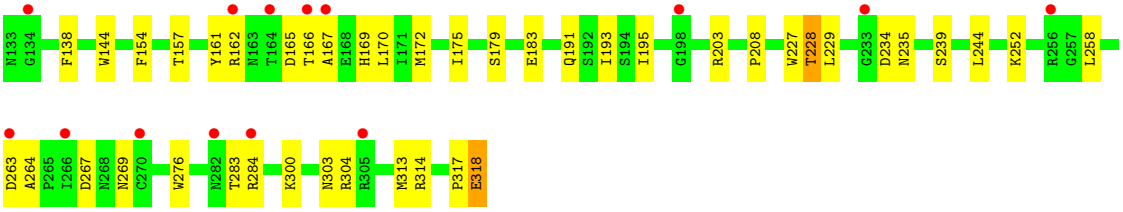


• Molecule 1: hemagglutinin

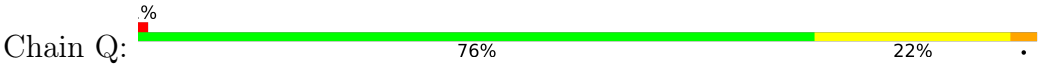


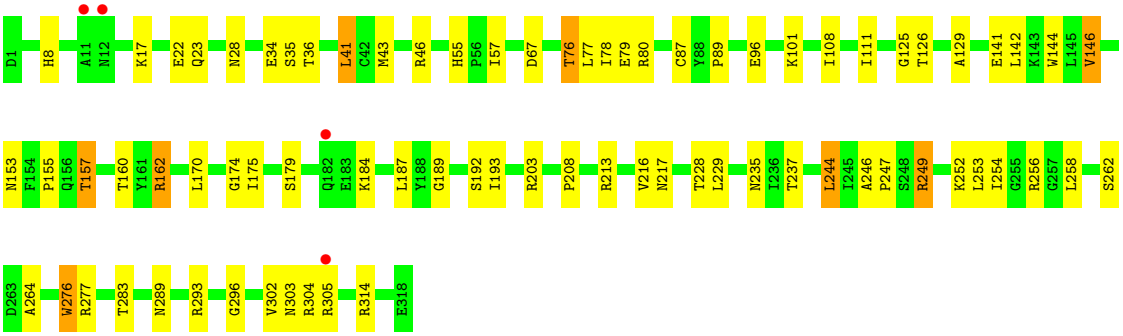
• Molecule 1: hemagglutinin



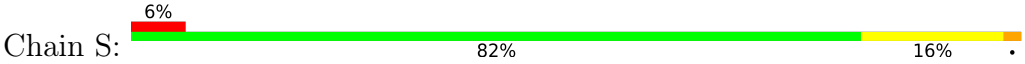


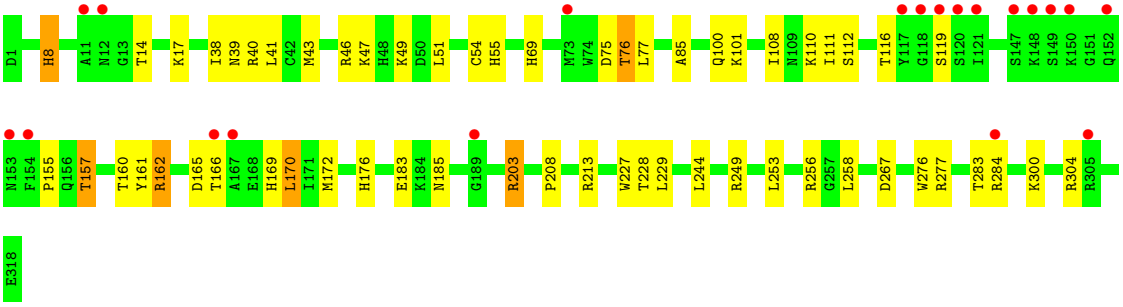
• Molecule 1: hemagglutinin



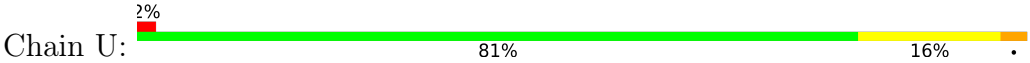


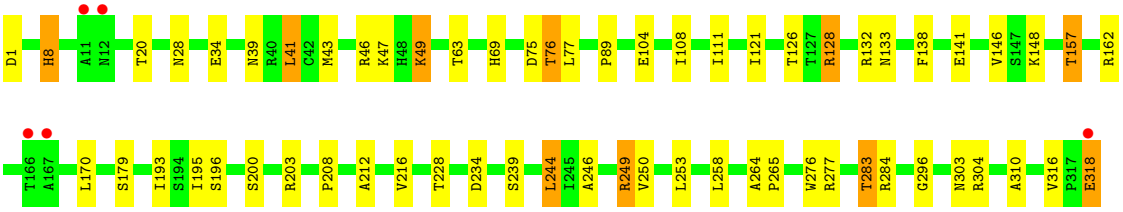
• Molecule 1: hemagglutinin



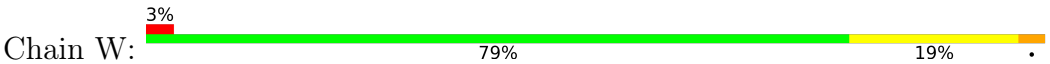


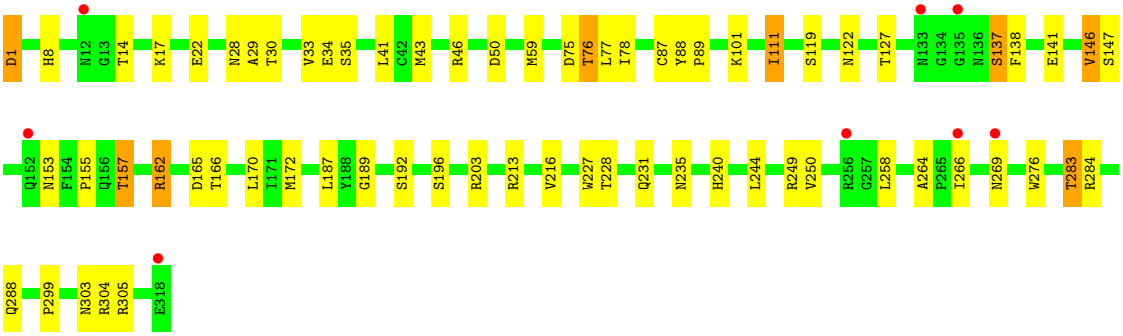
• Molecule 1: hemagglutinin



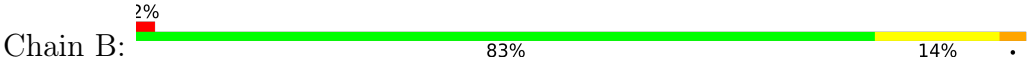


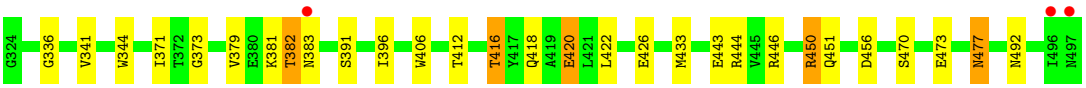
• Molecule 1: hemagglutinin



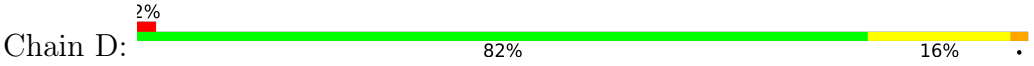


• Molecule 2: hemagglutinin

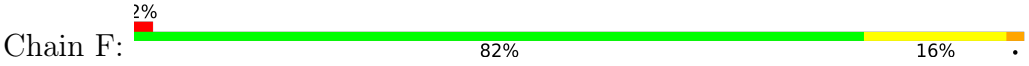


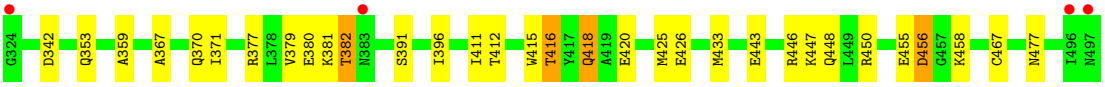


• Molecule 2: hemagglutinin

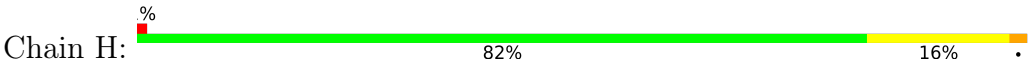


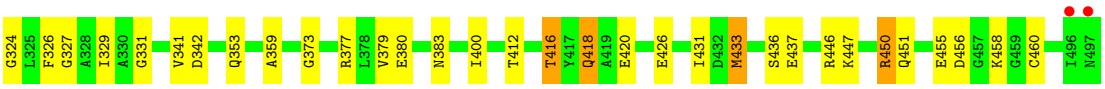
• Molecule 2: hemagglutinin



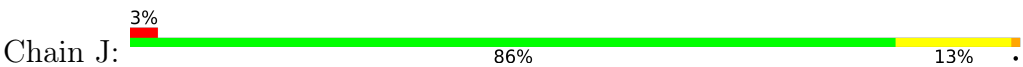


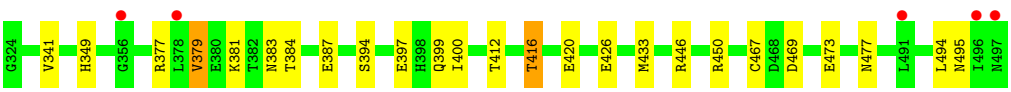
• Molecule 2: hemagglutinin





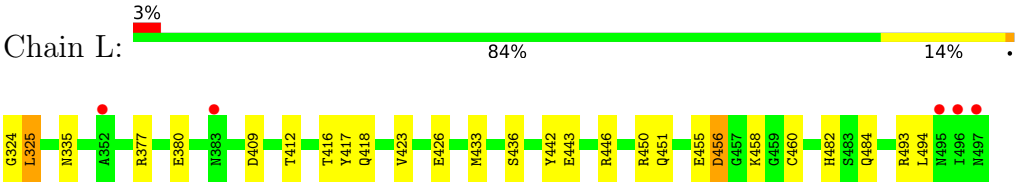
• Molecule 2: hemagglutinin





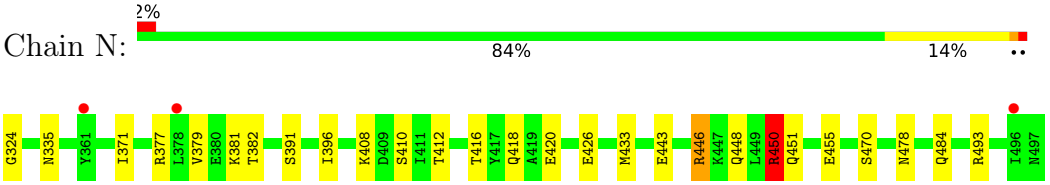
• Molecule 2: hemagglutinin

Chain L:



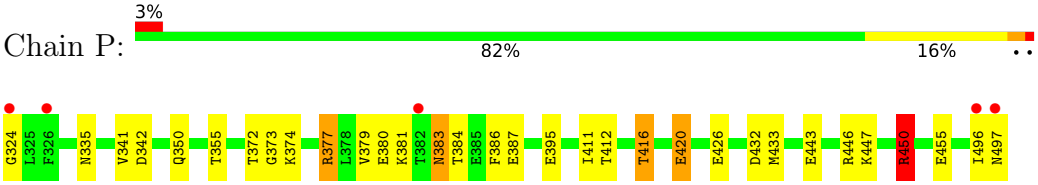
• Molecule 2: hemagglutinin

Chain N:



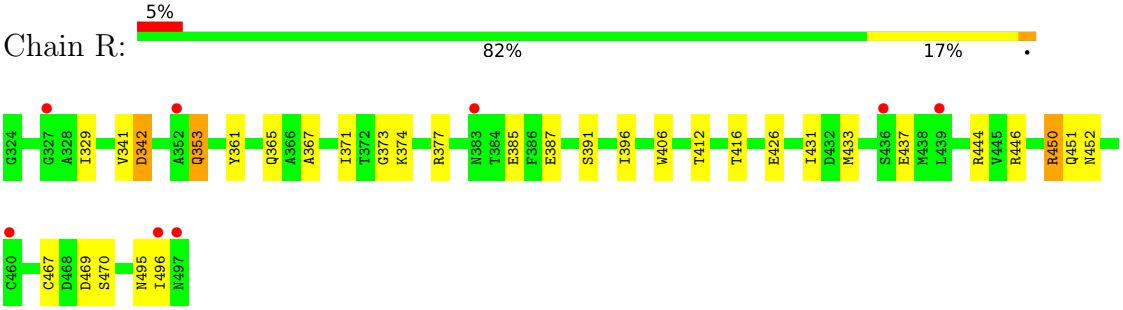
• Molecule 2: hemagglutinin

Chain P:



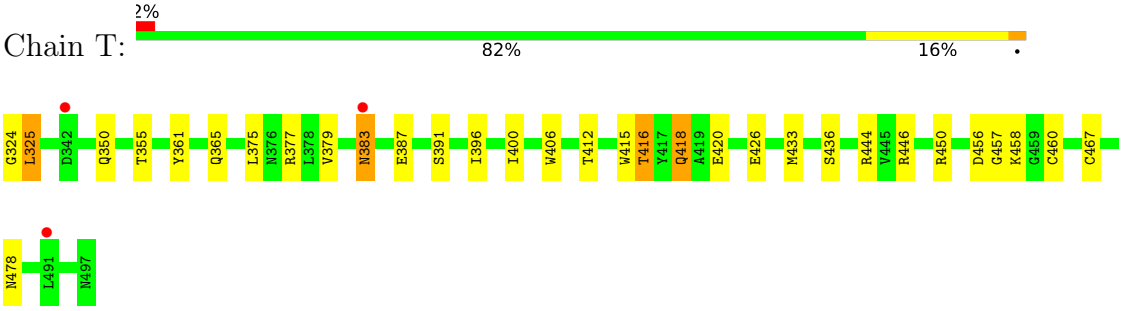
• Molecule 2: hemagglutinin

Chain R:



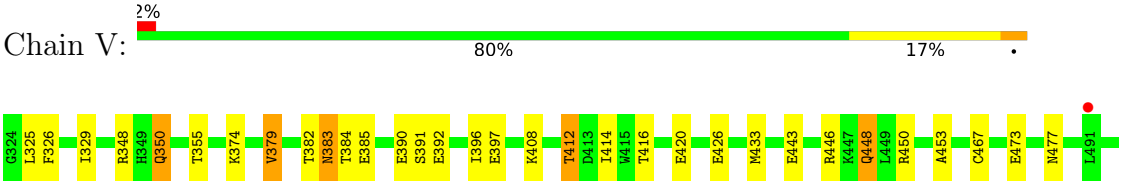
• Molecule 2: hemagglutinin

Chain T:



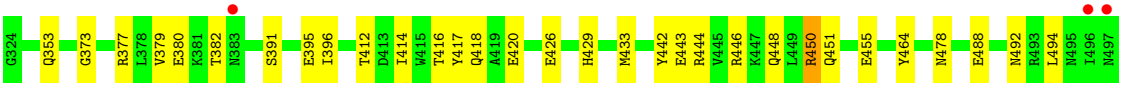
• Molecule 2: hemagglutinin

Chain V:



• Molecule 2: hemagglutinin

Chain X:



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

Chain Y:

• Molecule 4: N-acetyl-alpha-neuraminic acid-(2-3)-beta-D-galactopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain Z:

• Molecule 4: N-acetyl-alpha-neuraminic acid-(2-3)-beta-D-galactopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

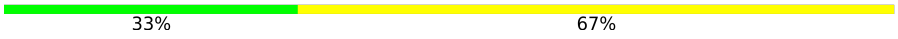
Chain a:

• Molecule 4: N-acetyl-alpha-neuraminic acid-(2-3)-beta-D-galactopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain b:

• Molecule 4: N-acetyl-alpha-neuraminic acid-(2-3)-beta-D-galactopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain c:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α, β, γ	93.26Å 111.44Å 222.35Å 94.96° 101.17° 91.90°	Depositor
Resolution (Å)	37.07 – 2.59 37.07 – 2.59	Depositor EDS
% Data completeness (in resolution range)	97.2 (37.07-2.59) 89.1 (37.07-2.59)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	2.03 (at 2.61Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.8.2_1309)	Depositor
R, R_{free}	0.197 , 0.246 0.197 , 0.247	Depositor DCC
R_{free} test set	13242 reflections (5.04%)	wwPDB-VP
Wilson B-factor (Å^2)	37.5	Xtriage
Anisotropy	0.604	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.33 , 43.2	EDS
L-test for twinning[2]	$< \|L\| > = 0.49$, $< L^2 > = 0.32$	Xtriage
Estimated twinning fraction	0.009 for h,-k,-h-l	Xtriage
F_o,F_c correlation	0.94	EDS
Total number of atoms	48138	wwPDB-VP
Average B, all atoms (Å^2)	38.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.62% of the height of the origin peak. No significant pseudotranslation is detected.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ 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: NAG, GAL, SIA

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	0.33	0/2486	0.50	0/3368
1	C	0.30	0/2486	0.48	0/3368
1	E	0.31	0/2486	0.49	0/3368
1	G	0.31	0/2486	0.48	0/3368
1	I	0.31	0/2486	0.49	0/3368
1	K	0.31	0/2486	0.49	0/3368
1	M	0.27	0/2486	0.45	0/3368
1	O	0.28	0/2486	0.46	0/3368
1	Q	0.31	0/2486	0.47	0/3368
1	S	0.29	0/2485	0.49	0/3366
1	U	0.30	0/2486	0.48	0/3368
1	W	0.29	0/2486	0.47	0/3368
2	B	0.33	0/1427	0.46	0/1926
2	D	0.34	0/1427	0.50	0/1926
2	F	0.34	0/1427	0.48	0/1926
2	H	0.33	0/1427	0.46	0/1926
2	J	0.31	0/1427	0.47	0/1926
2	L	0.32	0/1427	0.46	0/1926
2	N	0.31	0/1427	0.45	0/1926
2	P	0.33	0/1427	0.47	0/1926
2	R	0.31	0/1427	0.45	0/1926
2	T	0.32	0/1427	0.45	0/1926
2	V	0.31	0/1427	0.48	0/1926
2	X	0.30	0/1427	0.45	0/1926
All	All	0.31	0/46955	0.47	0/63526

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 (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	2437	0	2389	41	0
1	C	2437	0	2389	31	0
1	E	2437	0	2387	40	0
1	G	2437	0	2389	33	0
1	I	2437	0	2389	33	0
1	K	2437	0	2389	27	0
1	M	2437	0	2389	26	0
1	O	2437	0	2389	39	0
1	Q	2437	0	2389	38	0
1	S	2436	0	2389	33	0
1	U	2437	0	2389	34	0
1	W	2437	0	2389	37	0
2	B	1402	0	1298	21	0
2	D	1402	0	1298	27	0
2	F	1402	0	1298	20	0
2	H	1402	0	1298	29	0
2	J	1402	0	1298	20	0
2	L	1402	0	1298	21	0
2	N	1402	0	1298	19	0
2	P	1402	0	1298	22	0
2	R	1402	0	1298	26	0
2	T	1402	0	1298	23	0
2	V	1402	0	1298	26	0
2	X	1402	0	1298	23	0
3	Y	28	0	25	1	0
4	Z	45	0	38	4	0
4	a	45	0	38	0	0
4	b	45	0	38	0	0
4	c	45	0	38	0	0
5	A	14	0	13	0	0
5	B	14	0	13	0	0
5	C	14	0	13	0	0
5	D	14	0	13	1	0
5	E	28	0	26	1	0
5	F	14	0	13	0	0
5	G	14	0	13	0	0
5	H	14	0	13	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
5	I	14	0	13	0	0
5	J	14	0	13	1	0
5	K	14	0	13	1	0
5	L	14	0	13	0	0
5	M	14	0	13	0	0
5	N	14	0	13	1	0
5	O	14	0	13	1	0
5	P	14	0	13	0	0
5	Q	14	0	13	0	0
5	R	14	0	13	0	0
5	S	14	0	13	0	0
5	T	14	0	13	1	0
5	U	14	0	13	1	0
5	V	14	0	13	0	0
5	W	14	0	13	0	0
5	X	14	0	13	0	0
6	A	106	0	0	8	0
6	B	70	0	0	2	0
6	C	71	0	0	5	0
6	D	87	0	0	9	0
6	E	93	0	0	11	0
6	F	64	0	0	3	0
6	G	80	0	0	11	0
6	H	53	0	0	6	0
6	I	102	0	0	8	0
6	J	43	0	0	8	0
6	K	93	0	0	3	0
6	L	53	0	0	3	0
6	M	27	0	0	5	0
6	N	46	0	0	4	0
6	O	40	0	0	5	0
6	P	48	0	0	6	0
6	Q	84	0	0	5	0
6	R	42	0	0	8	0
6	S	64	0	0	7	0
6	T	49	0	0	3	0
6	U	76	0	0	7	0
6	V	44	0	0	9	0
6	W	50	0	0	11	0
6	X	28	0	0	4	0
All	All	48138	0	44744	609	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 7.

The worst 5 of 609 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)
2:L:409:ASP:OD2	6:L:738:HOH:O	1.81	0.98
1:A:96:GLU:OE1	6:A:712:HOH:O	1.86	0.92
2:R:444:ARG:O	6:R:723:HOH:O	1.89	0.89
2:V:390:GLU:OE2	6:V:709:HOH:O	1.90	0.89
1:G:11:ALA:O	6:G:712:HOH:O	1.91	0.88

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	316/318 (99%)	309 (98%)	7 (2%)	0	100	100
1	C	316/318 (99%)	306 (97%)	10 (3%)	0	100	100
1	E	316/318 (99%)	306 (97%)	10 (3%)	0	100	100
1	G	316/318 (99%)	310 (98%)	6 (2%)	0	100	100
1	I	316/318 (99%)	307 (97%)	9 (3%)	0	100	100
1	K	316/318 (99%)	306 (97%)	10 (3%)	0	100	100
1	M	316/318 (99%)	302 (96%)	13 (4%)	1 (0%)	41	64
1	O	316/318 (99%)	303 (96%)	13 (4%)	0	100	100
1	Q	316/318 (99%)	304 (96%)	12 (4%)	0	100	100
1	S	316/318 (99%)	305 (96%)	10 (3%)	1 (0%)	41	64
1	U	316/318 (99%)	307 (97%)	8 (2%)	1 (0%)	41	64
1	W	316/318 (99%)	305 (96%)	11 (4%)	0	100	100
2	B	172/174 (99%)	167 (97%)	4 (2%)	1 (1%)	25	47

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
2	D	172/174 (99%)	166 (96%)	4 (2%)	2 (1%)	13	27
2	F	172/174 (99%)	167 (97%)	4 (2%)	1 (1%)	25	47
2	H	172/174 (99%)	168 (98%)	3 (2%)	1 (1%)	25	47
2	J	172/174 (99%)	164 (95%)	7 (4%)	1 (1%)	25	47
2	L	172/174 (99%)	169 (98%)	2 (1%)	1 (1%)	25	47
2	N	172/174 (99%)	167 (97%)	4 (2%)	1 (1%)	25	47
2	P	172/174 (99%)	163 (95%)	8 (5%)	1 (1%)	25	47
2	R	172/174 (99%)	166 (96%)	5 (3%)	1 (1%)	25	47
2	T	172/174 (99%)	166 (96%)	5 (3%)	1 (1%)	25	47
2	V	172/174 (99%)	163 (95%)	8 (5%)	1 (1%)	25	47
2	X	172/174 (99%)	166 (96%)	5 (3%)	1 (1%)	25	47
All	All	5856/5904 (99%)	5662 (97%)	178 (3%)	16 (0%)	41	64

5 of 16 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	D	450	ARG
2	F	450	ARG
2	H	450	ARG
2	J	450	ARG
2	L	450	ARG

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	269/269 (100%)	248 (92%)	21 (8%)	12	25
1	C	269/269 (100%)	249 (93%)	20 (7%)	13	28
1	E	269/269 (100%)	247 (92%)	22 (8%)	11	22
1	G	269/269 (100%)	248 (92%)	21 (8%)	12	25
1	I	269/269 (100%)	250 (93%)	19 (7%)	14	29

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
1	K	269/269 (100%)	246 (91%)	23 (9%)	10 20
1	M	269/269 (100%)	250 (93%)	19 (7%)	14 29
1	O	269/269 (100%)	249 (93%)	20 (7%)	13 28
1	Q	269/269 (100%)	246 (91%)	23 (9%)	10 20
1	S	269/269 (100%)	250 (93%)	19 (7%)	14 29
1	U	269/269 (100%)	241 (90%)	28 (10%)	7 13
1	W	269/269 (100%)	245 (91%)	24 (9%)	9 19
2	B	148/148 (100%)	139 (94%)	9 (6%)	18 38
2	D	148/148 (100%)	138 (93%)	10 (7%)	16 32
2	F	148/148 (100%)	137 (93%)	11 (7%)	13 28
2	H	148/148 (100%)	144 (97%)	4 (3%)	44 71
2	J	148/148 (100%)	140 (95%)	8 (5%)	22 44
2	L	148/148 (100%)	140 (95%)	8 (5%)	22 44
2	N	148/148 (100%)	141 (95%)	7 (5%)	26 50
2	P	148/148 (100%)	137 (93%)	11 (7%)	13 28
2	R	148/148 (100%)	141 (95%)	7 (5%)	26 50
2	T	148/148 (100%)	141 (95%)	7 (5%)	26 50
2	V	148/148 (100%)	139 (94%)	9 (6%)	18 38
2	X	148/148 (100%)	143 (97%)	5 (3%)	37 63
All	All	5004/5004 (100%)	4649 (93%)	355 (7%)	14 29

5 of 355 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	P	433	MET
2	T	433	MET
1	Q	111	ILE
2	R	467	CYS
1	U	148	LYS

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

Mol	Chain	Res	Type
1	S	163	ASN

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Mol	Chain	Res	Type
2	T	418	GLN
2	V	448	GLN
1	U	303	ASN
2	F	370	GLN

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)

14 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	Y	1	3,1	14,14,15	0.24	0	17,19,21	0.44	0
3	NAG	Y	2	3	14,14,15	0.84	0	17,19,21	0.70	0
4	NAG	Z	1	4	14,14,15	0.31	0	17,19,21	0.39	0
4	GAL	Z	2	4	11,11,12	0.78	0	15,15,17	1.02	0
4	SIA	Z	3	4	20,20,21	0.58	0	24,28,31	0.76	0
4	NAG	a	1	4	14,14,15	0.26	0	17,19,21	0.35	0
4	GAL	a	2	4	11,11,12	0.28	0	15,15,17	0.77	0
4	SIA	a	3	4	20,20,21	0.59	0	24,28,31	0.76	0
4	NAG	b	1	4	14,14,15	0.38	0	17,19,21	0.47	0
4	GAL	b	2	4	11,11,12	0.91	0	15,15,17	1.30	2 (13%)
4	SIA	b	3	4	20,20,21	0.57	0	24,28,31	0.76	0
4	NAG	c	1	4	14,14,15	0.24	0	17,19,21	0.48	0
4	GAL	c	2	4	11,11,12	0.76	0	15,15,17	1.12	1 (6%)
4	SIA	c	3	4	20,20,21	1.98	9 (45%)	24,28,31	3.04	11 (45%)

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	Y	1	3,1	-	2/6/23/26	0/1/1/1
3	NAG	Y	2	3	-	1/6/23/26	0/1/1/1
4	NAG	Z	1	4	-	2/6/23/26	0/1/1/1
4	GAL	Z	2	4	-	2/2/19/22	0/1/1/1
4	SIA	Z	3	4	-	1/18/34/38	0/1/1/1
4	NAG	a	1	4	-	2/6/23/26	0/1/1/1
4	GAL	a	2	4	-	0/2/19/22	0/1/1/1
4	SIA	a	3	4	-	0/18/34/38	0/1/1/1
4	NAG	b	1	4	-	2/6/23/26	0/1/1/1
4	GAL	b	2	4	-	2/2/19/22	0/1/1/1
4	SIA	b	3	4	-	0/18/34/38	0/1/1/1
4	NAG	c	1	4	-	2/6/23/26	0/1/1/1
4	GAL	c	2	4	-	0/2/19/22	0/1/1/1
4	SIA	c	3	4	-	6/18/34/38	0/1/1/1

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

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	c	3	SIA	C8-C7	-3.48	1.46	1.53
4	c	3	SIA	C4-C5	-3.08	1.50	1.53
4	c	3	SIA	O10-C10	-3.03	1.16	1.23
4	c	3	SIA	C7-C6	-2.71	1.49	1.53
4	c	3	SIA	C3-C4	2.59	1.57	1.52

The worst 5 of 14 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
4	c	3	SIA	O6-C2-C3	-9.09	97.95	110.46
4	c	3	SIA	C6-O6-C2	5.52	123.15	111.34
4	c	3	SIA	O9-C9-C8	-5.34	99.45	111.07
4	c	3	SIA	O10-C10-C11	-4.01	114.62	122.06
4	c	3	SIA	C11-C10-N5	3.86	122.63	116.10

There are no chirality outliers.

5 of 22 torsion outliers are listed below:

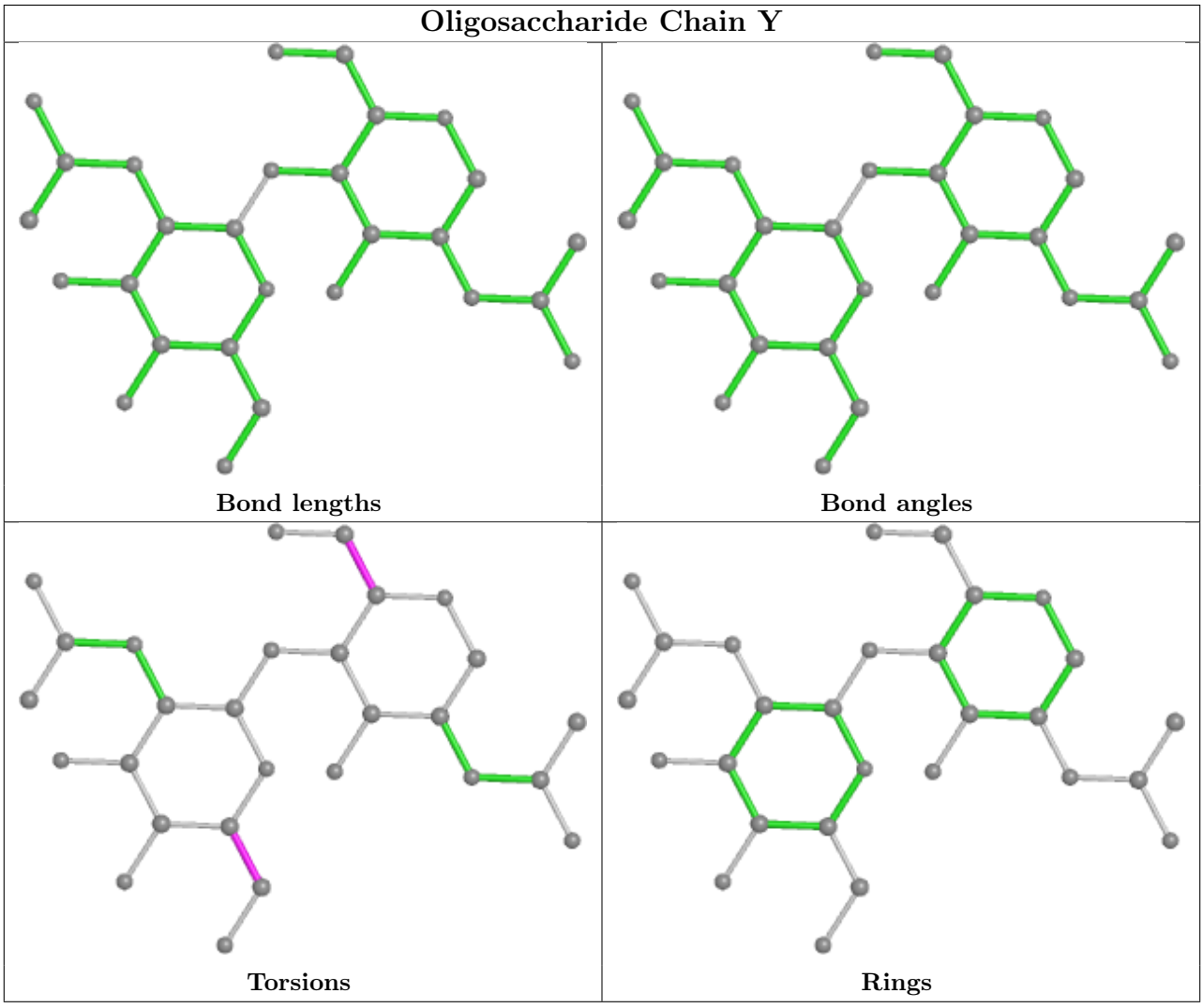
Mol	Chain	Res	Type	Atoms
4	c	3	SIA	O6-C6-C7-C8
4	c	3	SIA	O6-C6-C7-O7
3	Y	1	NAG	O5-C5-C6-O6
4	b	1	NAG	O5-C5-C6-O6
4	b	2	GAL	O5-C5-C6-O6

There are no ring outliers.

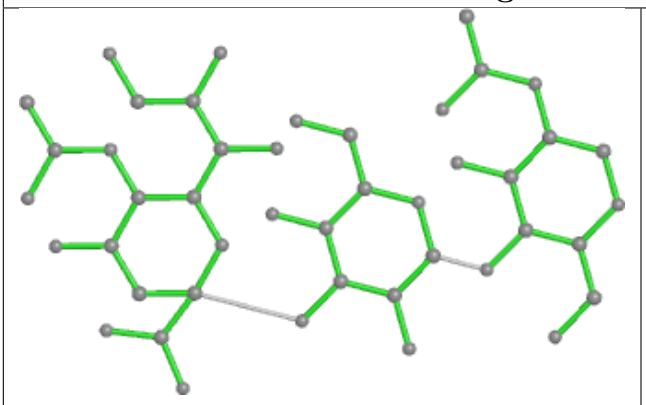
2 monomers are involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	Z	3	SIA	4	0
3	Y	1	NAG	1	0

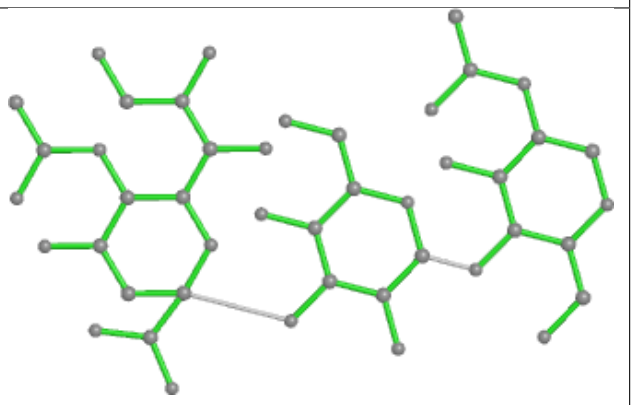
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for oligosaccharide.



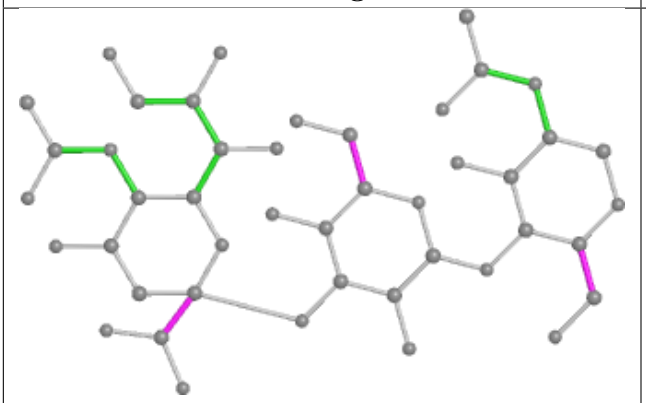
Oligosaccharide Chain Z



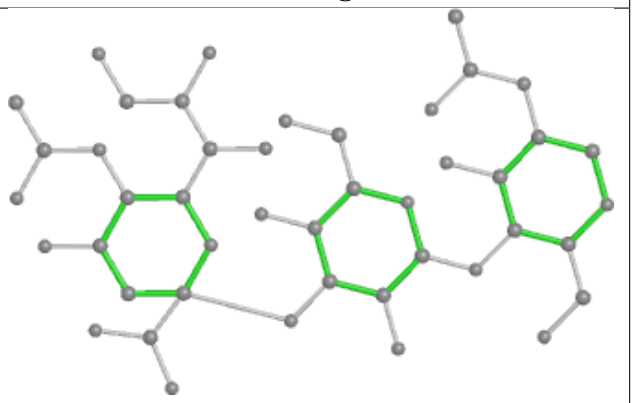
Bond lengths



Bond angles

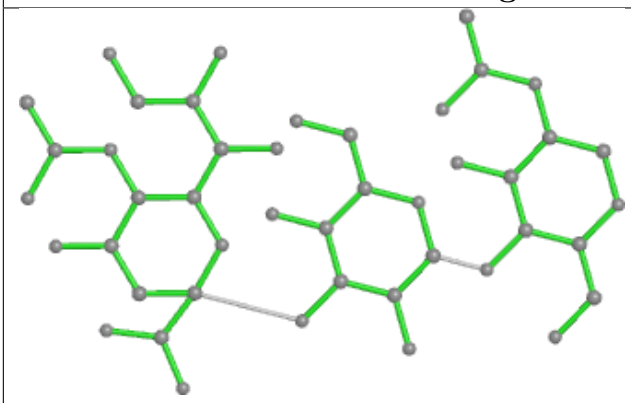


Torsions

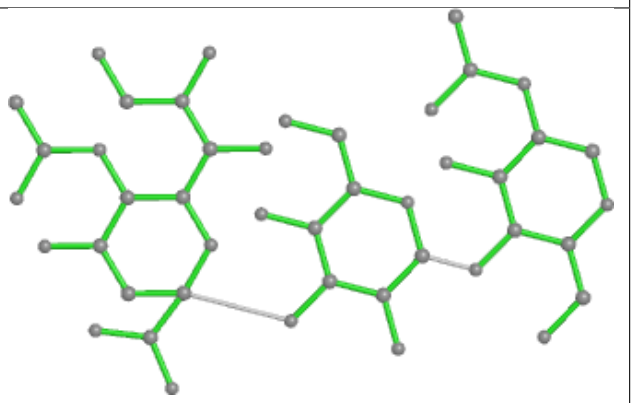


Rings

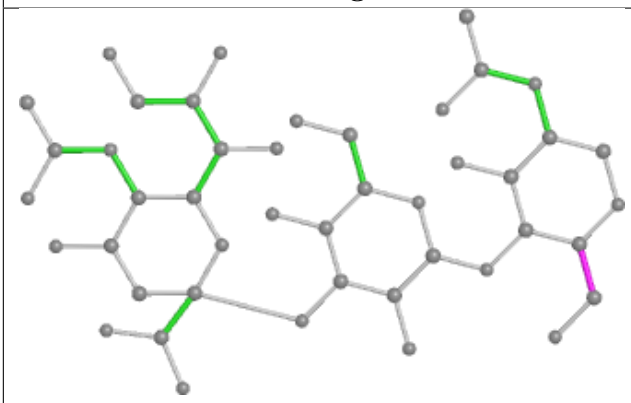
Oligosaccharide Chain a



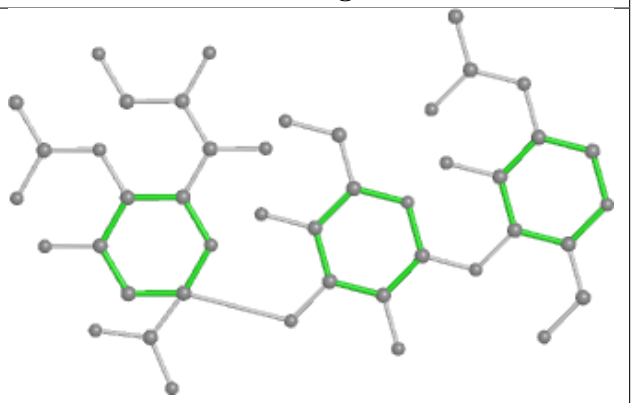
Bond lengths



Bond angles

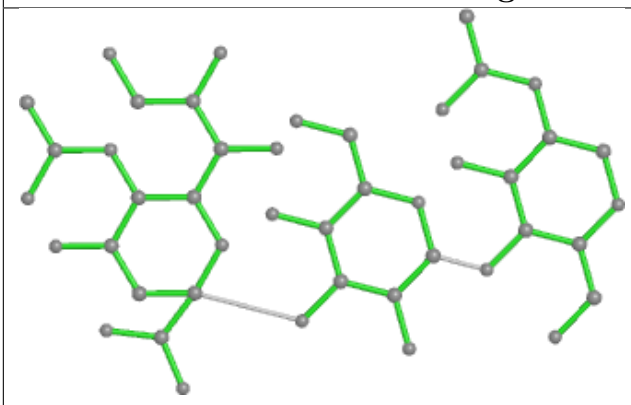


Torsions

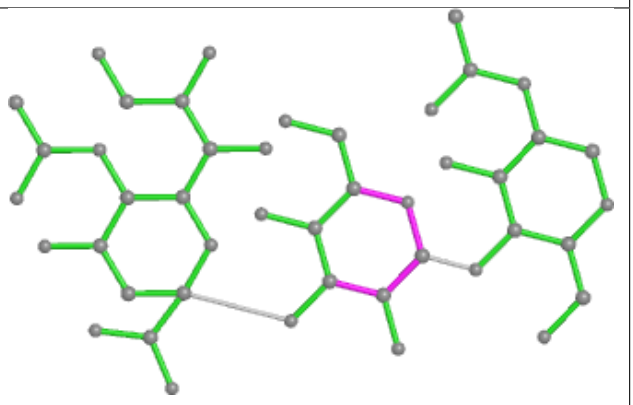


Rings

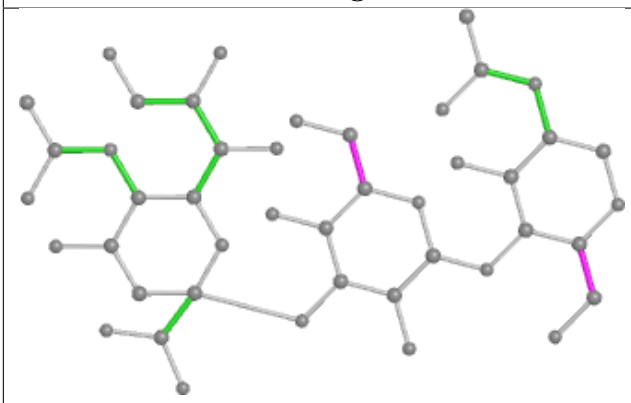
Oligosaccharide Chain b



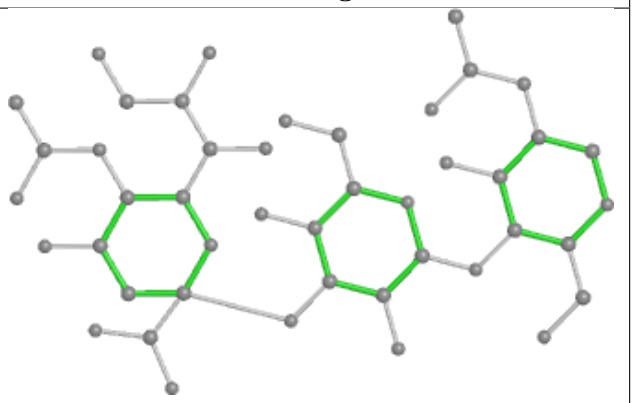
Bond lengths



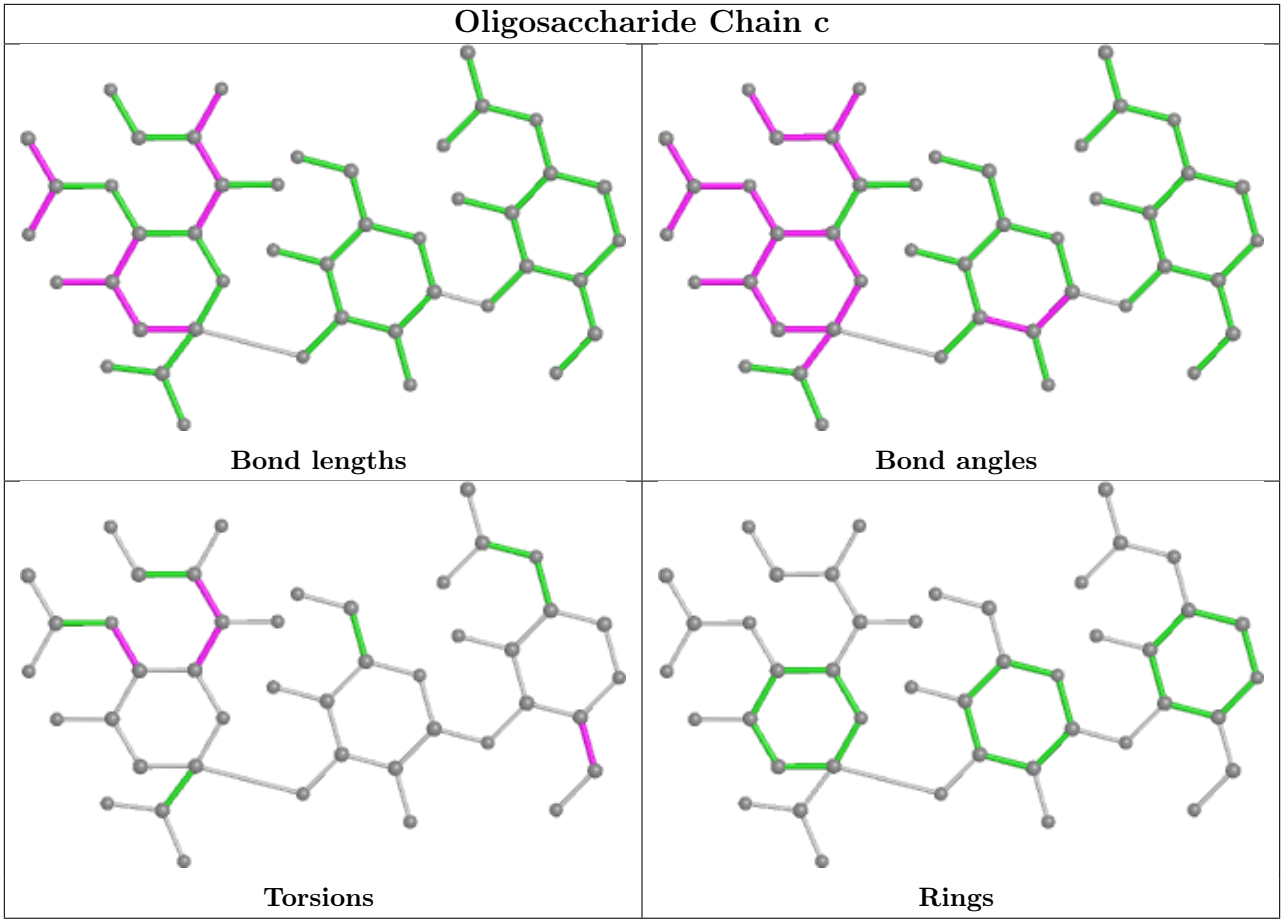
Bond angles



Torsions



Rings



5.6 Ligand geometry (i)

25 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	$\#\|Z\| > 2$
5	NAG	L	601	2	14,14,15	0.37	0	17,19,21	0.47	0
5	NAG	B	601	2	14,14,15	0.35	0	17,19,21	0.51	0
5	NAG	I	601	1	14,14,15	0.23	0	17,19,21	0.45	0
5	NAG	N	601	2	14,14,15	0.47	0	17,19,21	0.44	0
5	NAG	W	601	1	14,14,15	0.29	0	17,19,21	0.61	0
5	NAG	M	601	1	14,14,15	0.36	0	17,19,21	0.57	0
5	NAG	U	601	1	14,14,15	0.42	0	17,19,21	0.43	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\|>2$	Counts	RMSZ	$\#\|Z\|>2$
5	NAG	H	601	2	14,14,15	0.42	0	17,19,21	0.59	0
5	NAG	Q	601	1	14,14,15	0.28	0	17,19,21	0.40	0
5	NAG	C	601	1	14,14,15	0.21	0	17,19,21	0.41	0
5	NAG	G	601	1	14,14,15	0.18	0	17,19,21	0.46	0
5	NAG	P	601	2	14,14,15	0.36	0	17,19,21	0.48	0
5	NAG	R	601	2	14,14,15	0.21	0	17,19,21	0.62	0
5	NAG	O	601	1	14,14,15	0.20	0	17,19,21	0.37	0
5	NAG	K	601	1	14,14,15	0.35	0	17,19,21	0.54	0
5	NAG	D	601	2	14,14,15	0.30	0	17,19,21	0.44	0
5	NAG	F	601	2	14,14,15	0.31	0	17,19,21	0.61	0
5	NAG	X	601	2	14,14,15	0.51	0	17,19,21	0.54	0
5	NAG	J	601	2	14,14,15	0.70	0	17,19,21	0.76	0
5	NAG	V	601	2	14,14,15	0.40	0	17,19,21	0.51	0
5	NAG	S	601	1	14,14,15	0.28	0	17,19,21	0.55	0
5	NAG	A	601	1	14,14,15	0.29	0	17,19,21	0.53	0
5	NAG	E	601	1	14,14,15	0.38	0	17,19,21	0.50	0
5	NAG	T	601	2	14,14,15	0.55	0	17,19,21	0.40	0
5	NAG	E	602	1	14,14,15	0.90	1 (7%)	17,19,21	0.93	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
5	NAG	L	601	2	-	2/6/23/26	0/1/1/1
5	NAG	B	601	2	-	0/6/23/26	0/1/1/1
5	NAG	I	601	1	-	0/6/23/26	0/1/1/1
5	NAG	N	601	2	-	0/6/23/26	0/1/1/1
5	NAG	W	601	1	-	4/6/23/26	0/1/1/1
5	NAG	M	601	1	-	0/6/23/26	0/1/1/1
5	NAG	U	601	1	-	2/6/23/26	0/1/1/1
5	NAG	H	601	2	-	0/6/23/26	0/1/1/1
5	NAG	Q	601	1	-	2/6/23/26	0/1/1/1
5	NAG	C	601	1	-	0/6/23/26	0/1/1/1
5	NAG	G	601	1	-	2/6/23/26	0/1/1/1
5	NAG	P	601	2	-	0/6/23/26	0/1/1/1
5	NAG	R	601	2	-	0/6/23/26	0/1/1/1
5	NAG	O	601	1	-	2/6/23/26	0/1/1/1
5	NAG	K	601	1	-	2/6/23/26	0/1/1/1
5	NAG	D	601	2	-	0/6/23/26	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
5	NAG	F	601	2	-	0/6/23/26	0/1/1/1
5	NAG	X	601	2	-	2/6/23/26	0/1/1/1
5	NAG	J	601	2	-	0/6/23/26	0/1/1/1
5	NAG	V	601	2	-	0/6/23/26	0/1/1/1
5	NAG	S	601	1	-	0/6/23/26	0/1/1/1
5	NAG	A	601	1	-	2/6/23/26	0/1/1/1
5	NAG	E	601	1	-	2/6/23/26	0/1/1/1
5	NAG	T	601	2	-	2/6/23/26	0/1/1/1
5	NAG	E	602	1	-	0/6/23/26	0/1/1/1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	E	602	NAG	C2-N2	2.26	1.50	1.46

There are no bond angle outliers.

There are no chirality outliers.

5 of 24 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	W	601	NAG	C8-C7-N2-C2
5	W	601	NAG	O7-C7-N2-C2
5	O	601	NAG	O5-C5-C6-O6
5	T	601	NAG	O5-C5-C6-O6
5	X	601	NAG	O5-C5-C6-O6

There are no ring outliers.

8 monomers are involved in 8 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	N	601	NAG	1	0
5	U	601	NAG	1	0
5	O	601	NAG	1	0
5	K	601	NAG	1	0
5	D	601	NAG	1	0
5	J	601	NAG	1	0
5	E	601	NAG	1	0
5	T	601	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

6.1 Protein, DNA and RNA chains

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^{th} 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	318/318 (100%)	-0.19	0 100 100	10, 25, 42, 69	0
1	C	318/318 (100%)	0.11	4 (1%) 77 73	10, 36, 64, 95	0
1	E	318/318 (100%)	-0.05	1 (0%) 94 93	11, 30, 48, 70	0
1	G	318/318 (100%)	0.04	4 (1%) 77 73	15, 34, 54, 80	0
1	I	318/318 (100%)	-0.03	5 (1%) 72 68	11, 29, 52, 101	0
1	K	318/318 (100%)	-0.08	6 (1%) 66 62	13, 29, 52, 84	0
1	M	318/318 (100%)	0.63	31 (9%) 7 5	24, 56, 87, 123	0
1	O	318/318 (100%)	0.48	25 (7%) 12 9	22, 52, 82, 112	0
1	Q	318/318 (100%)	-0.10	4 (1%) 77 73	12, 33, 54, 92	0
1	S	318/318 (100%)	0.31	20 (6%) 20 15	17, 43, 70, 107	0
1	U	318/318 (100%)	-0.05	5 (1%) 72 68	16, 31, 53, 83	0
1	W	318/318 (100%)	0.22	8 (2%) 57 51	24, 44, 71, 94	0
2	B	174/174 (100%)	0.23	3 (1%) 70 66	12, 26, 54, 108	0
2	D	174/174 (100%)	0.24	3 (1%) 70 66	11, 25, 52, 130	0
2	F	174/174 (100%)	0.17	4 (2%) 60 54	10, 27, 56, 101	0
2	H	174/174 (100%)	0.18	2 (1%) 80 78	12, 29, 55, 125	0
2	J	174/174 (100%)	0.40	5 (2%) 51 45	12, 37, 74, 132	0
2	L	174/174 (100%)	0.31	5 (2%) 51 45	14, 35, 67, 108	0
2	N	174/174 (100%)	0.31	3 (1%) 70 66	23, 38, 71, 109	0
2	P	174/174 (100%)	0.36	5 (2%) 51 45	18, 37, 66, 125	0
2	R	174/174 (100%)	0.29	8 (4%) 32 26	16, 33, 55, 110	0
2	T	174/174 (100%)	0.19	3 (1%) 70 66	13, 37, 61, 119	0
2	V	174/174 (100%)	0.29	3 (1%) 70 66	14, 39, 71, 127	0
2	X	174/174 (100%)	0.35	3 (1%) 70 66	19, 45, 69, 119	0

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Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å^2)	Q<0.9
All	All	5904/5904 (100%)	0.17	160 (2%) 54 48	10, 35, 69, 132	0

The worst 5 of 160 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	D	497	ASN	13.2
2	P	497	ASN	12.7
2	J	497	ASN	11.3
2	V	496	ILE	10.5
2	X	497	ASN	10.3

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)

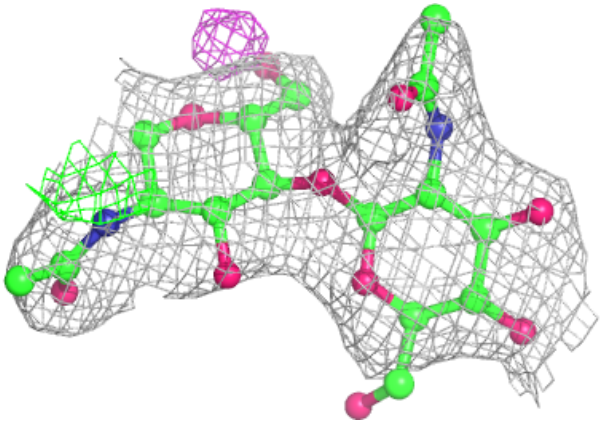
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

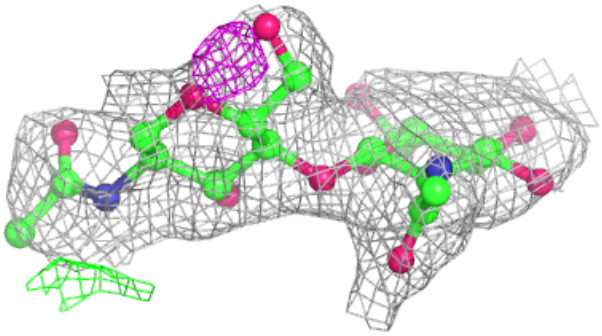
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
4	NAG	a	1	14/15	0.70	0.78	125,140,147,147	0
4	NAG	Z	1	14/15	0.72	0.56	100,109,111,112	0
4	GAL	a	2	11/12	0.74	0.24	72,90,112,114	0
4	NAG	c	1	14/15	0.76	0.58	99,115,125,128	0
4	NAG	b	1	14/15	0.77	0.49	97,103,110,110	0
4	GAL	b	2	11/12	0.78	0.24	66,76,99,100	0
4	GAL	c	2	11/12	0.78	0.27	58,69,81,85	0
4	GAL	Z	2	11/12	0.86	0.24	90,98,116,120	0
4	SIA	Z	3	20/21	0.87	0.19	34,45,58,60	0
4	SIA	a	3	20/21	0.87	0.24	44,57,71,72	0
3	NAG	Y	2	14/15	0.89	0.26	48,69,77,80	0
4	SIA	b	3	20/21	0.90	0.19	26,35,46,47	0
4	SIA	c	3	20/21	0.93	0.15	24,34,46,47	0
3	NAG	Y	1	14/15	0.94	0.22	46,54,64,68	0

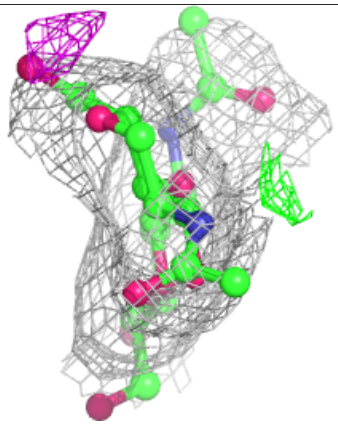
The following is a graphical depiction of the model fit to experimental electron density for oligosaccharide. Each fit is shown from different orientation to approximate a three-dimensional view.

Electron density around Chain Y:

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

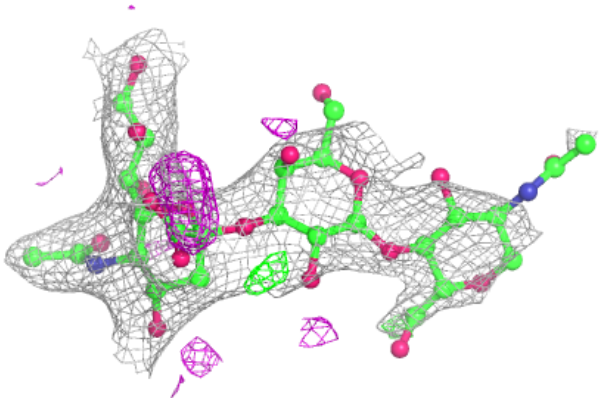


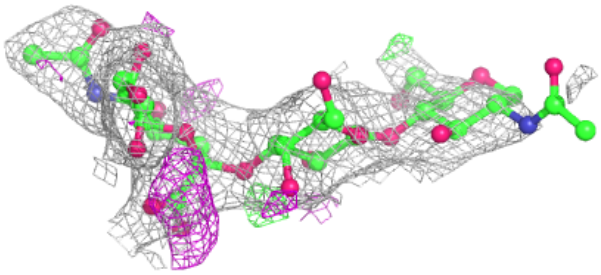


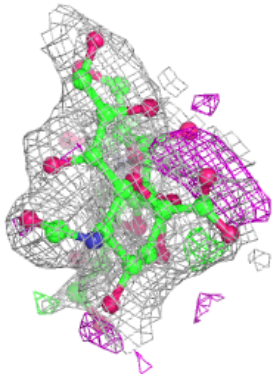


Electron density around Chain Z:

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

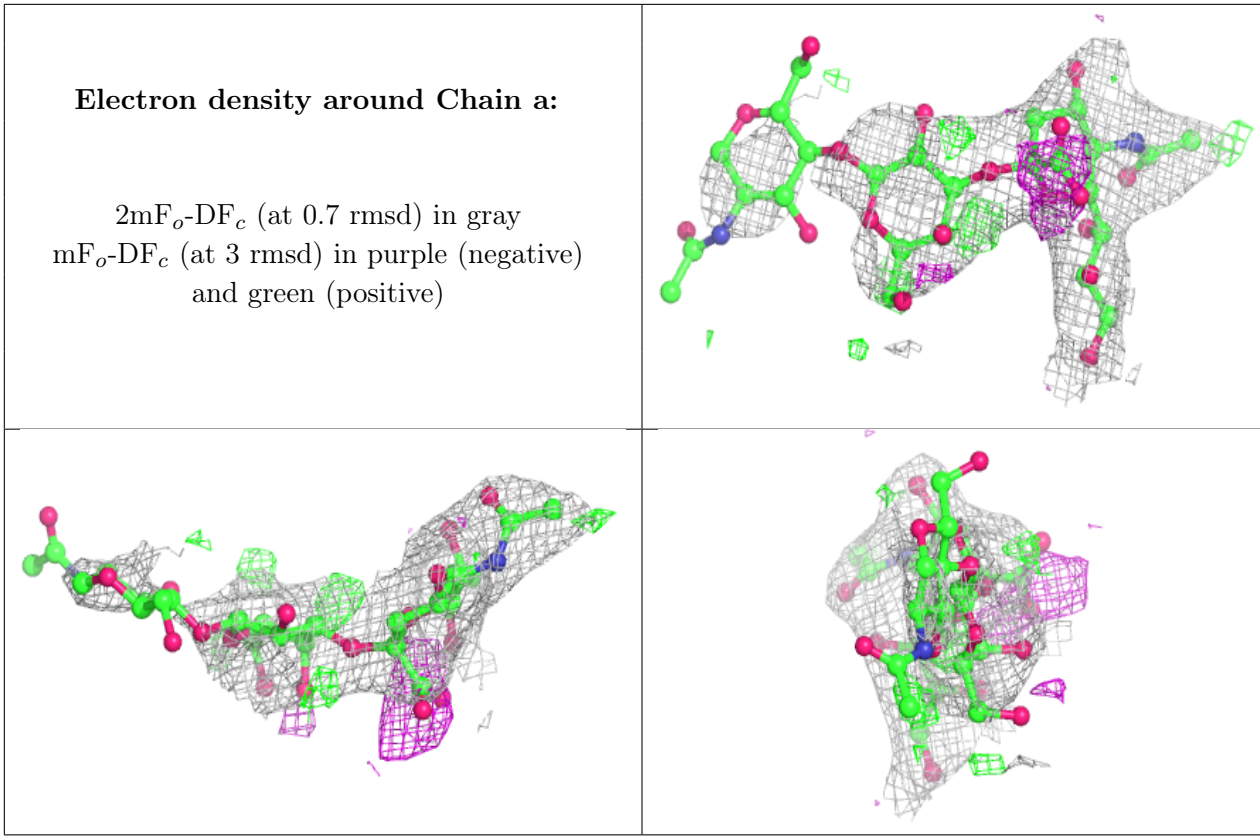






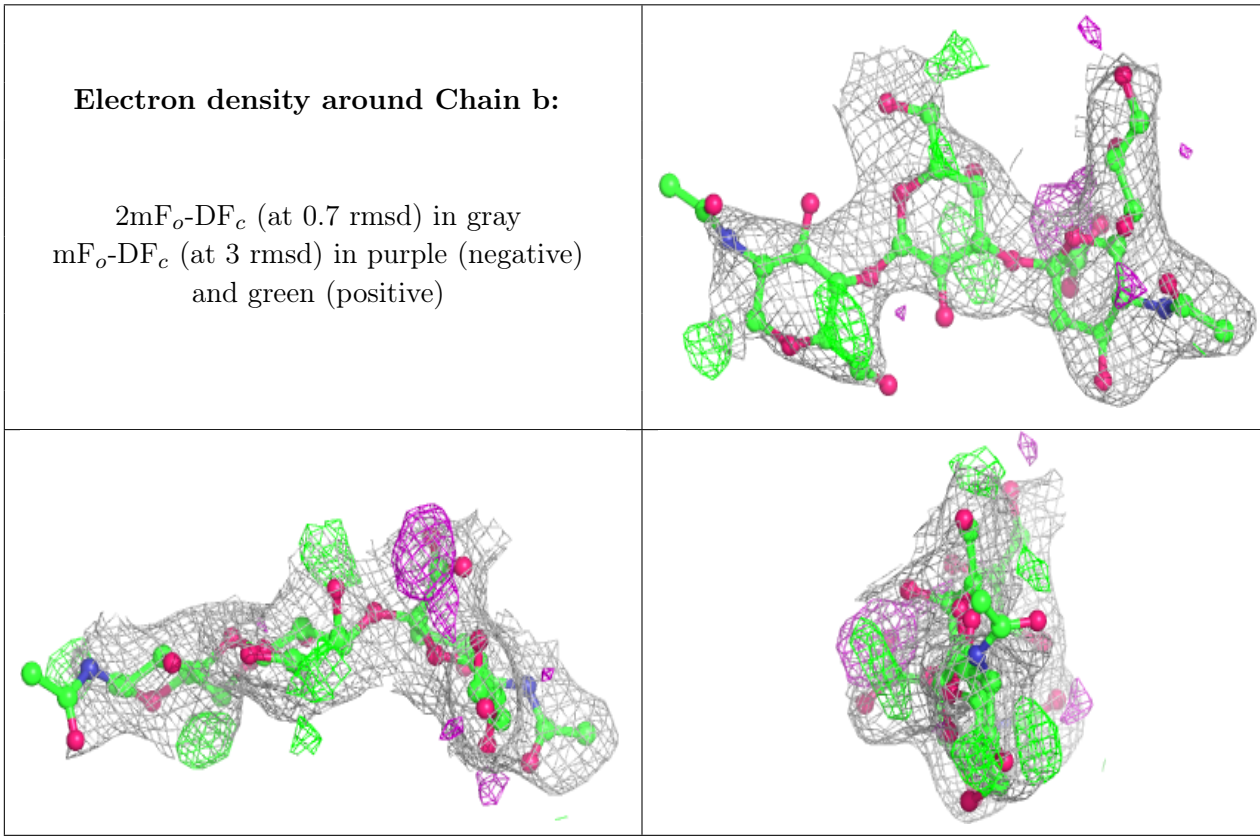
Electron density around Chain a:

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



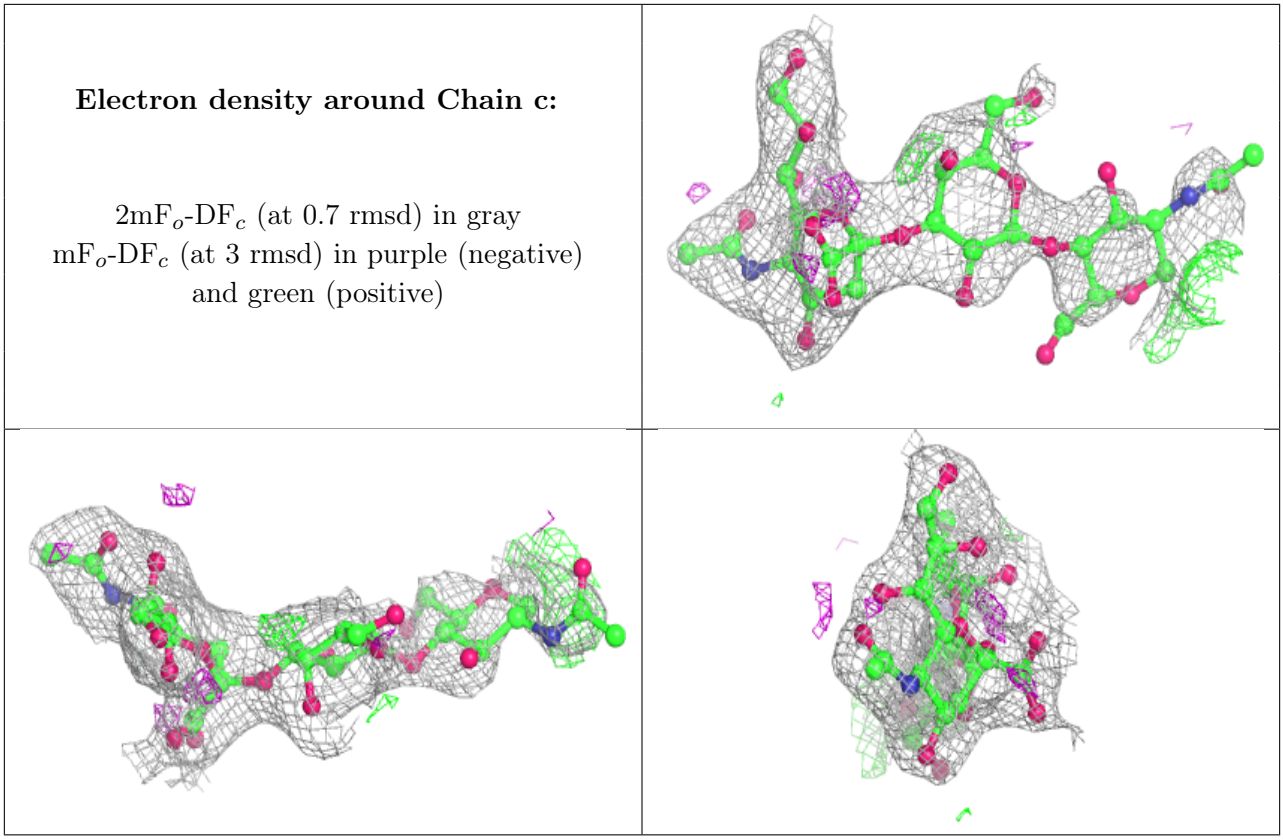
Electron density around Chain b:

$2mF_o$-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 c:

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



6.4 Ligands

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^{th} 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
5	NAG	O	601	14/15	0.72	0.30	61,88,97,101	0
5	NAG	W	601	14/15	0.72	0.30	79,87,91,92	0
5	NAG	S	601	14/15	0.79	0.20	54,75,88,92	0
5	NAG	U	601	14/15	0.84	0.32	80,98,105,113	0
5	NAG	L	601	14/15	0.85	0.17	23,32,38,38	0
5	NAG	R	601	14/15	0.86	0.15	32,39,49,50	0
5	NAG	G	601	14/15	0.86	0.17	43,64,72,74	0
5	NAG	C	601	14/15	0.87	0.19	46,67,77,77	0
5	NAG	M	601	14/15	0.87	0.28	66,83,89,92	0
5	NAG	K	601	14/15	0.87	0.26	46,71,77,79	0
5	NAG	Q	601	14/15	0.87	0.20	76,87,98,99	0
5	NAG	X	601	14/15	0.87	0.16	37,46,54,54	0
5	NAG	F	601	14/15	0.88	0.14	25,30,33,36	0
5	NAG	E	602	14/15	0.88	0.27	54,58,73,77	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
5	NAG	N	601	14/15	0.90	0.13	40,50,51,54	0
5	NAG	V	601	14/15	0.90	0.17	28,38,40,44	0
5	NAG	E	601	14/15	0.91	0.14	48,62,69,70	0
5	NAG	H	601	14/15	0.91	0.13	22,30,33,33	0
5	NAG	I	601	14/15	0.92	0.17	52,62,65,65	0
5	NAG	A	601	14/15	0.93	0.18	34,57,68,74	0
5	NAG	B	601	14/15	0.93	0.10	29,38,43,50	0
5	NAG	P	601	14/15	0.93	0.13	34,38,44,44	0
5	NAG	T	601	14/15	0.93	0.12	23,34,41,43	0
5	NAG	D	601	14/15	0.95	0.11	20,28,36,37	0
5	NAG	J	601	14/15	0.95	0.13	18,28,36,38	0

6.5 Other polymers (i)

There are no such residues in this entry.