wwPDB X-ray Structure Validation Summary Report (i)

Sep 5, 2023 – 04:37 AM EDT

PDB ID	:	3UBQ
Title	:	Influenza hemagglutinin from the 2009 pandemic in complex with ligand 3SLN
Authors	:	Xu, R.; Wilson, I.A.
Deposited on	:	2011-10-24
Resolution	:	2.00 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:

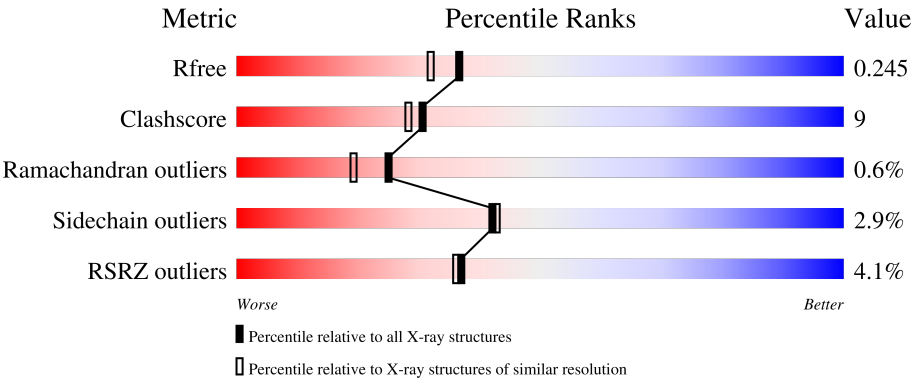
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.35
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.35

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 2.00 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.



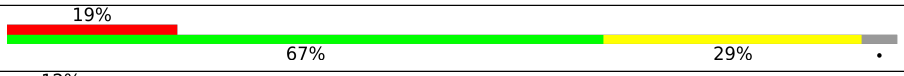
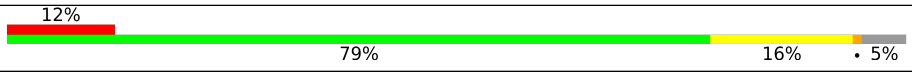
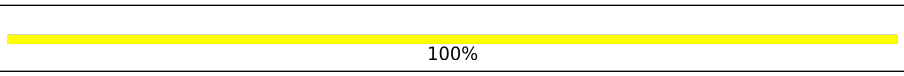
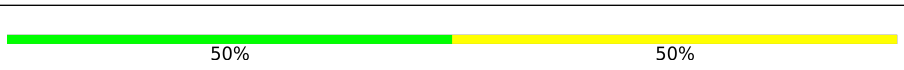
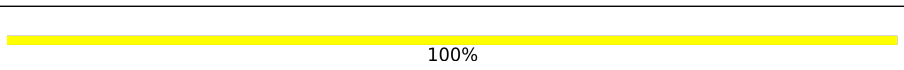
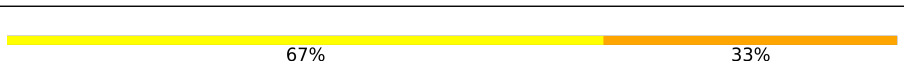
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	8085 (2.00-2.00)
Clashscore	141614	9178 (2.00-2.00)
Ramachandran outliers	138981	9054 (2.00-2.00)
Sidechain outliers	138945	9053 (2.00-2.00)
RSRZ outliers	127900	7900 (2.00-2.00)

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	329	2% 75% 21%
1	C	329	2% 74% 22%
1	E	329	% 76% 20%
1	G	329	% 78% 19%
1	I	329	% 81% 15%

Continued on next page...

Continued from previous page...

Mol	Chain	Length	Quality of chain
1	K	329	
2	B	177	
2	D	177	
2	F	177	
2	H	177	
2	J	177	
2	L	177	
3	M	2	
3	O	2	
3	R	2	
4	N	3	
5	P	2	
5	Q	2	

2 Entry composition (i)

There are 7 unique types of molecules in this entry. The entry contains 24675 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.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	323	Total	C	N	O	S	0	0	0
			2522	1595	432	482	13			
1	C	321	Total	C	N	O	S	0	0	0
			2509	1588	430	478	13			
1	E	321	Total	C	N	O	S	0	0	0
			2504	1585	430	476	13			
1	G	323	Total	C	N	O	S	0	0	0
			2522	1595	432	482	13			
1	I	319	Total	C	N	O	S	0	0	0
			2495	1579	428	475	13			
1	K	318	Total	C	N	O	S	0	0	0
			2490	1576	427	474	13			

There are 24 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	9	PRO	-	expression tag	UNP C3W5S1
A	10	GLY	-	expression tag	UNP C3W5S1
A	205	CYS	GLY	engineered mutation	UNP C3W5S1
A	220	CYS	ARG	engineered mutation	UNP C3W5S1
C	9	PRO	-	expression tag	UNP C3W5S1
C	10	GLY	-	expression tag	UNP C3W5S1
C	205	CYS	GLY	engineered mutation	UNP C3W5S1
C	220	CYS	ARG	engineered mutation	UNP C3W5S1
E	9	PRO	-	expression tag	UNP C3W5S1
E	10	GLY	-	expression tag	UNP C3W5S1
E	205	CYS	GLY	engineered mutation	UNP C3W5S1
E	220	CYS	ARG	engineered mutation	UNP C3W5S1
G	9	PRO	-	expression tag	UNP C3W5S1
G	10	GLY	-	expression tag	UNP C3W5S1
G	205	CYS	GLY	engineered mutation	UNP C3W5S1
G	220	CYS	ARG	engineered mutation	UNP C3W5S1
I	9	PRO	-	expression tag	UNP C3W5S1

Continued on next page...

Continued from previous page...

Chain	Residue	Modelled	Actual	Comment	Reference
I	10	GLY	-	expression tag	UNP C3W5S1
I	205	CYS	GLY	engineered mutation	UNP C3W5S1
I	220	CYS	ARG	engineered mutation	UNP C3W5S1
K	9	PRO	-	expression tag	UNP C3W5S1
K	10	GLY	-	expression tag	UNP C3W5S1
K	205	CYS	GLY	engineered mutation	UNP C3W5S1
K	220	CYS	ARG	engineered mutation	UNP C3W5S1

- Molecule 2 is a protein called hemagglutinin HA2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	169	Total	C	N	O	S	0	0	0
			1360	855	229	270	6			
2	D	170	Total	C	N	O	S	0	0	0
			1371	861	233	271	6			
2	F	170	Total	C	N	O	S	0	0	0
			1371	861	233	271	6			
2	H	172	Total	C	N	O	S	0	0	0
			1389	871	235	277	6			
2	J	170	Total	C	N	O	S	0	0	0
			1371	861	233	271	6			
2	L	169	Total	C	N	O	S	0	0	0
			1360	855	229	270	6			

There are 18 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	175	SER	-	expression tag	UNP C3W5S1
B	176	GLY	-	expression tag	UNP C3W5S1
B	177	ARG	-	expression tag	UNP C3W5S1
D	175	SER	-	expression tag	UNP C3W5S1
D	176	GLY	-	expression tag	UNP C3W5S1
D	177	ARG	-	expression tag	UNP C3W5S1
F	175	SER	-	expression tag	UNP C3W5S1
F	176	GLY	-	expression tag	UNP C3W5S1
F	177	ARG	-	expression tag	UNP C3W5S1
H	175	SER	-	expression tag	UNP C3W5S1
H	176	GLY	-	expression tag	UNP C3W5S1
H	177	ARG	-	expression tag	UNP C3W5S1
J	175	SER	-	expression tag	UNP C3W5S1
J	176	GLY	-	expression tag	UNP C3W5S1
J	177	ARG	-	expression tag	UNP C3W5S1

Continued on next page...

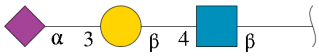
Continued from previous page...

Chain	Residue	Modelled	Actual	Comment	Reference
L	175	SER	-	expression tag	UNP C3W5S1
L	176	GLY	-	expression tag	UNP C3W5S1
L	177	ARG	-	expression tag	UNP C3W5S1

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

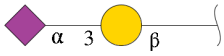
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
3	M	2	Total	C	N	O	0	0	0
			28	16	2	10			
3	O	2	Total	C	N	O	0	0	0
			28	16	2	10			
3	R	2	Total	C	N	O	0	0	0
			28	16	2	10			

- 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	N	3	Total	C	N	O	0	0	0
			46	25	2	19			

- Molecule 5 is an oligosaccharide called N-acetyl-alpha-neuraminic acid-(2-3)-beta-D-galactopyranose.



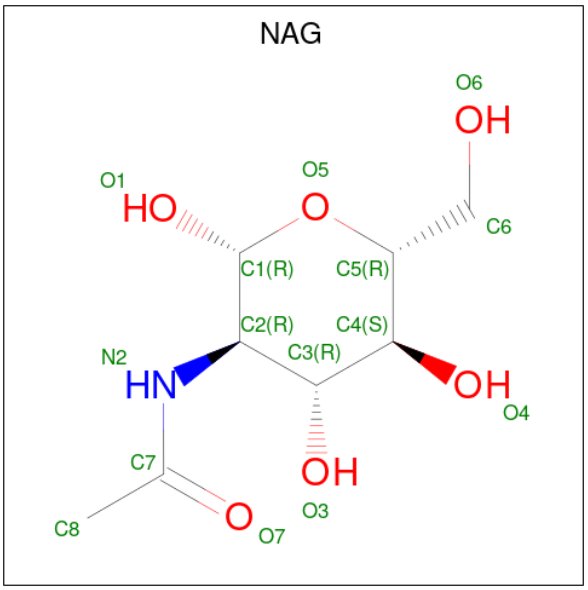
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
5	P	2	Total	C	N	O	0	0	0
			31	17	1	13			

Continued on next page...

Continued from previous page...

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf	Trace
5	Q	2	Total C N O 31 17 1 13	0	0	0

- 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 14 8 1 5	0	0
6	D	1	Total C N O 14 8 1 5	0	0
6	E	1	Total C N O 14 8 1 5	0	0
6	G	1	Total C N O 14 8 1 5	0	0
6	G	1	Total C N O 14 8 1 5	0	0
6	G	1	Total C N O 14 8 1 5	0	0
6	I	1	Total C N O 14 8 1 5	0	0
6	I	1	Total C N O 14 8 1 5	0	0
6	I	1	Total C N O 14 8 1 5	0	0
6	K	1	Total C N O 14 8 1 5	0	0

Continued on next page...

Continued from previous page...

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

- Molecule 7 is water.

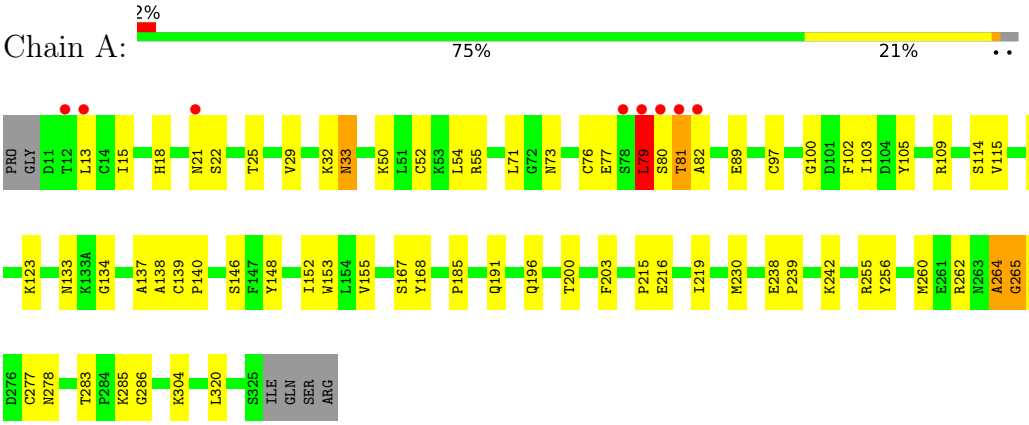
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
7	A	138	Total O 138 138	0	0
7	B	54	Total O 54 54	0	0
7	C	89	Total O 89 89	0	0
7	D	66	Total O 66 66	0	0
7	E	117	Total O 117 117	0	0
7	F	62	Total O 62 62	0	0
7	G	121	Total O 121 121	0	0
7	H	69	Total O 69 69	0	0
7	I	127	Total O 127 127	0	0
7	J	57	Total O 57 57	0	0
7	K	97	Total O 97 97	0	0
7	L	54	Total O 54 54	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.

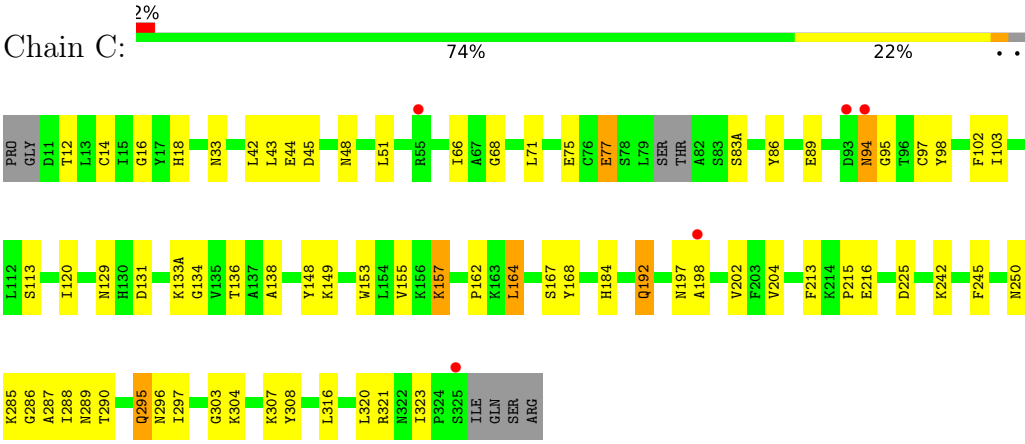
• Molecule 1: hemagglutinin HA1

Chain A:



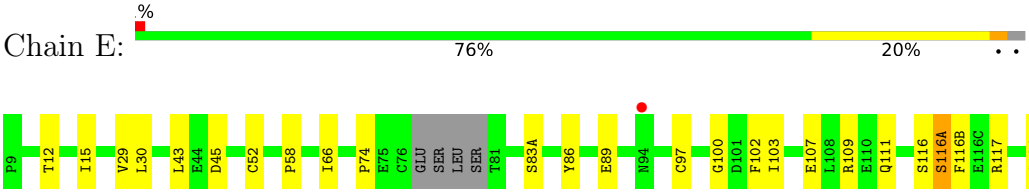
• Molecule 1: hemagglutinin HA1

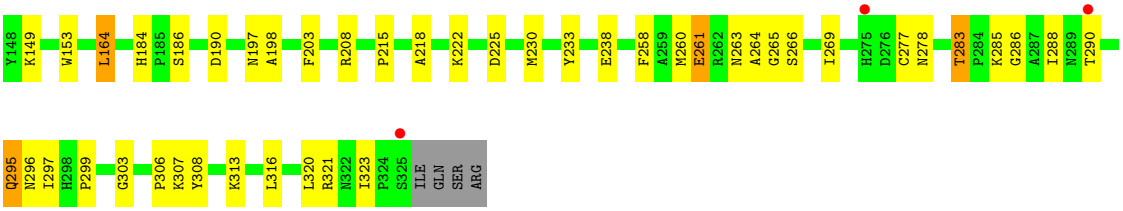
Chain C:



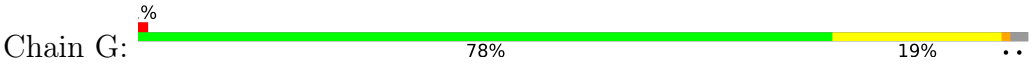
• Molecule 1: hemagglutinin HA1

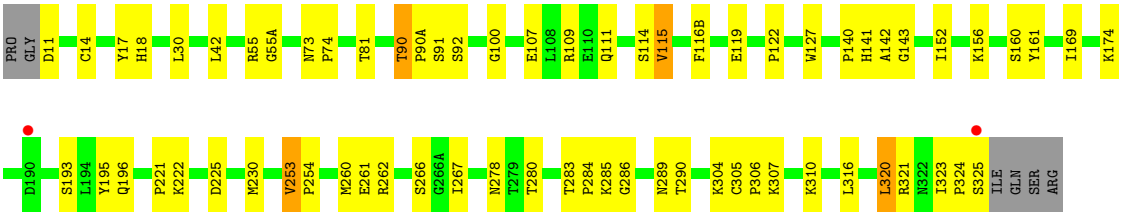
Chain E:



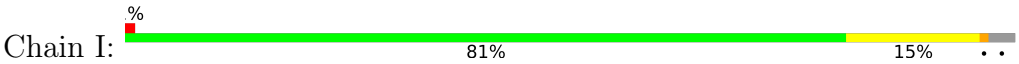


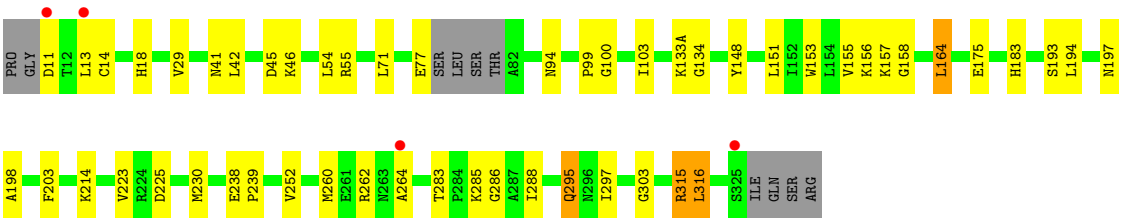
• Molecule 1: hemagglutinin HA1



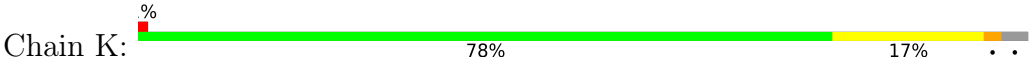


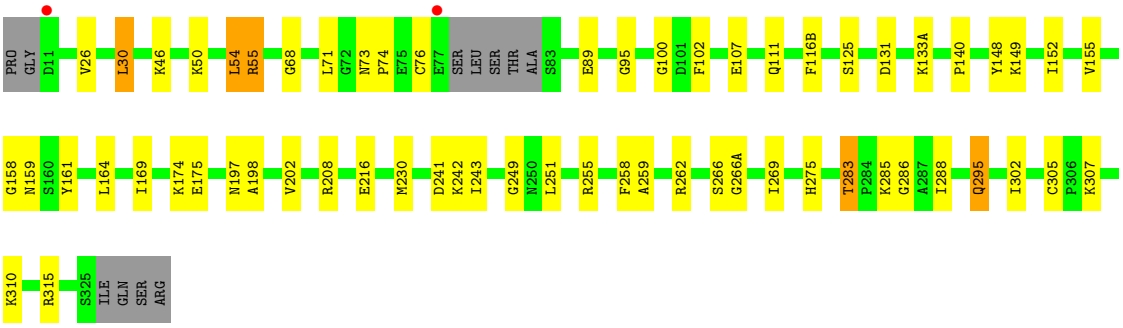
• Molecule 1: hemagglutinin HA1



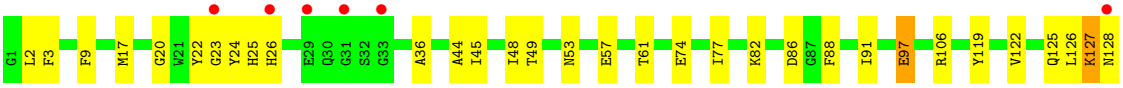


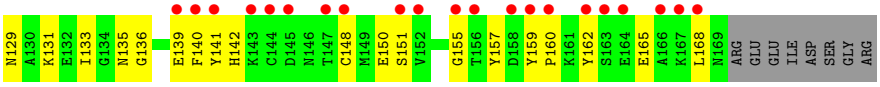
• Molecule 1: hemagglutinin HA1





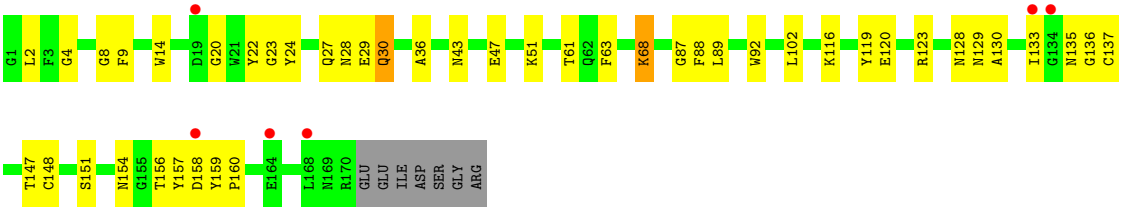
• Molecule 2: hemagglutinin HA2





• Molecule 2: hemagglutinin HA2

Chain D:

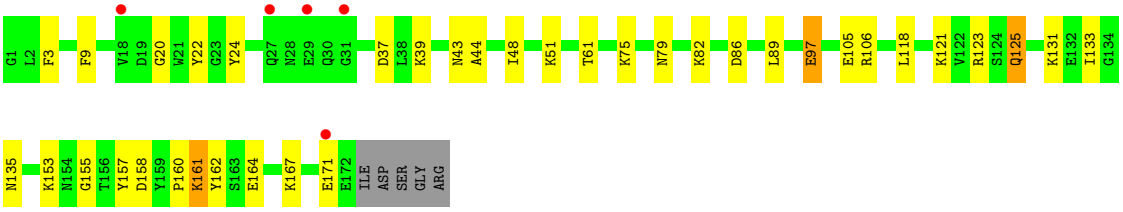


• Molecule 2: hemagglutinin HA2

Chain F:

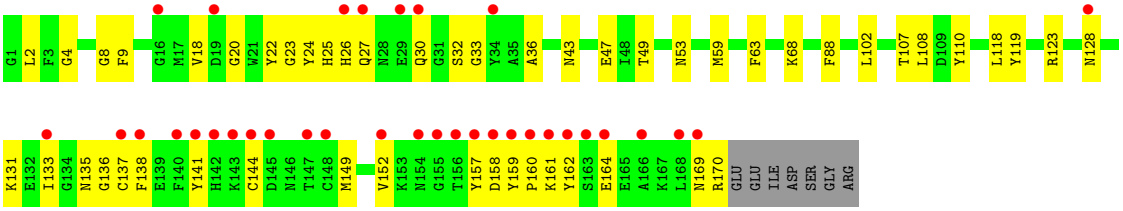
• Molecule 2: hemagglutinin HA2

Chain H:



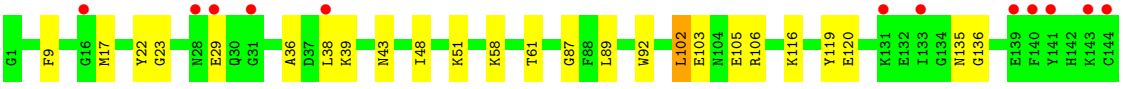
• Molecule 2: hemagglutinin HA2

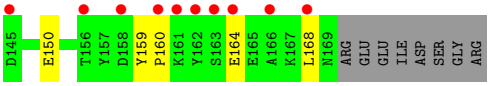
Chain J:



• Molecule 2: hemagglutinin HA2

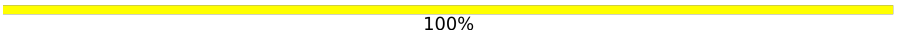
Chain L:





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

Chain M:

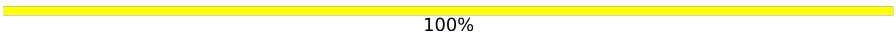


• Molecule 3: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain O:

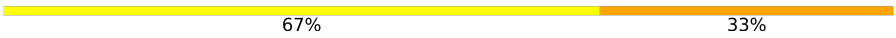
• Molecule 3: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain R:



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

Chain N:



• Molecule 5: N-acetyl-alpha-neuraminic acid-(2-3)-beta-D-galactopyranose

Chain P:

• Molecule 5: N-acetyl-alpha-neuraminic acid-(2-3)-beta-D-galactopyranose

Chain Q:

4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α, β, γ	66.88Å 116.26Å 118.37Å 60.96° 77.13° 80.40°	Depositor
Resolution (Å)	49.24 – 2.00 49.24 – 2.00	Depositor EDS
% Data completeness (in resolution range)	84.2 (49.24-2.00) 84.2 (49.24-2.00)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	1.67 (at 2.00Å)	Xtriage
Refinement program	PHENIX 1.7_650	Depositor
R, R_{free}	0.199 , 0.249 0.197 , 0.245	Depositor DCC
R_{free} test set	8649 reflections (5.01%)	wwPDB-VP
Wilson B-factor (Å^2)	29.5	Xtriage
Anisotropy	0.168	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.35 , 48.2	EDS
L-test for twinning[2]	$< \|L\| > = 0.49$, $< L^2 > = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.95	EDS
Total number of atoms	24675	wwPDB-VP
Average B, all atoms (Å^2)	41.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.13% 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: SIA, GAL, 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	0.30	0/2586	0.49	0/3516
1	C	0.27	0/2572	0.47	0/3495
1	E	0.30	0/2568	0.51	0/3490
1	G	0.31	0/2586	0.49	0/3516
1	I	0.30	0/2558	0.49	0/3476
1	K	0.28	0/2553	0.48	0/3469
2	B	0.27	0/1388	0.42	0/1871
2	D	0.28	0/1399	0.43	0/1885
2	F	0.29	0/1399	0.43	0/1885
2	H	0.29	0/1417	0.43	0/1909
2	J	0.27	0/1399	0.42	0/1885
2	L	0.27	0/1388	0.41	0/1871
All	All	0.29	0/23813	0.47	0/32268

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2522	0	2466	61	0

Continued on next page...

Continued from previous page...

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	C	2509	0	2455	56	0
1	E	2504	0	2449	54	0
1	G	2522	0	2465	52	0
1	I	2495	0	2436	39	0
1	K	2490	0	2430	49	0
2	B	1360	0	1284	31	0
2	D	1371	0	1296	36	0
2	F	1371	0	1297	21	0
2	H	1389	0	1309	33	0
2	J	1371	0	1297	35	0
2	L	1360	0	1284	20	0
3	M	28	0	25	0	0
3	O	28	0	25	0	0
3	R	28	0	25	2	0
4	N	46	0	40	1	0
5	P	31	0	26	2	0
5	Q	31	0	26	1	0
6	A	14	0	13	1	0
6	D	14	0	13	0	0
6	E	14	0	13	0	0
6	G	42	0	39	3	0
6	I	42	0	39	0	0
6	K	42	0	39	0	0
7	A	138	0	0	5	0
7	B	54	0	0	1	0
7	C	89	0	0	2	0
7	D	66	0	0	1	0
7	E	117	0	0	3	0
7	F	62	0	0	0	0
7	G	121	0	0	2	0
7	H	69	0	0	2	0
7	I	127	0	0	1	0
7	J	57	0	0	0	0
7	K	97	0	0	1	0
7	L	54	0	0	2	0
All	All	24675	0	22791	426	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 426 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:H:97:GLU:HG2	2:L:58:LYS:HD2	1.37	1.06
1:C:283:THR:HG22	1:C:285:LYS:H	1.20	1.03
1:I:283:THR:HG22	1:I:285:LYS:H	1.20	1.02
1:A:283:THR:HG22	1:A:285:LYS:H	1.28	0.97
1:A:79:LEU:HD12	1:A:81:THR:HB	1.49	0.92

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	321/329 (98%)	303 (94%)	13 (4%)	5 (2%)	9 4
1	C	317/329 (96%)	301 (95%)	15 (5%)	1 (0%)	41 37
1	E	317/329 (96%)	306 (96%)	10 (3%)	1 (0%)	41 37
1	G	321/329 (98%)	305 (95%)	16 (5%)	0	100 100
1	I	315/329 (96%)	300 (95%)	12 (4%)	3 (1%)	15 9
1	K	314/329 (95%)	295 (94%)	17 (5%)	2 (1%)	25 19
2	B	167/177 (94%)	161 (96%)	5 (3%)	1 (1%)	25 19
2	D	168/177 (95%)	160 (95%)	6 (4%)	2 (1%)	13 7
2	F	168/177 (95%)	159 (95%)	9 (5%)	0	100 100
2	H	170/177 (96%)	164 (96%)	6 (4%)	0	100 100
2	J	168/177 (95%)	157 (94%)	9 (5%)	2 (1%)	13 7
2	L	167/177 (94%)	157 (94%)	10 (6%)	0	100 100
All	All	2913/3036 (96%)	2768 (95%)	128 (4%)	17 (1%)	25 19

5 of 17 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	81	THR

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	A	265	GLY
2	B	127	LYS
1	C	94	ASN
1	E	116(A)	SER

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	285/290 (98%)	280 (98%)	5 (2%)	59	63
1	C	283/290 (98%)	272 (96%)	11 (4%)	32	30
1	E	282/290 (97%)	273 (97%)	9 (3%)	39	38
1	G	285/290 (98%)	277 (97%)	8 (3%)	43	44
1	I	281/290 (97%)	274 (98%)	7 (2%)	47	49
1	K	281/290 (97%)	274 (98%)	7 (2%)	47	49
2	B	145/152 (95%)	139 (96%)	6 (4%)	30	28
2	D	146/152 (96%)	141 (97%)	5 (3%)	37	36
2	F	146/152 (96%)	144 (99%)	2 (1%)	67	72
2	H	148/152 (97%)	143 (97%)	5 (3%)	37	36
2	J	146/152 (96%)	140 (96%)	6 (4%)	30	28
2	L	145/152 (95%)	141 (97%)	4 (3%)	43	44
All	All	2573/2652 (97%)	2498 (97%)	75 (3%)	42	43

5 of 75 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	I	316	LEU
2	L	22	TYR
2	J	43	ASN
1	K	54	LEU
2	D	68	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 18 such sidechains are listed below:

Mol	Chain	Res	Type
2	J	53	ASN
2	L	95	ASN
2	L	60	ASN
1	G	196	GLN
2	J	27	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)

13 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	M	1	3,1	14,14,15	0.60	0	17,19,21	1.01	1 (5%)
3	NAG	M	2	3	14,14,15	0.55	0	17,19,21	0.77	1 (5%)
4	NAG	N	1	4	15,15,15	0.49	0	21,21,21	0.78	1 (4%)
4	GAL	N	2	4	11,11,12	0.68	0	15,15,17	1.22	2 (13%)
4	SIA	N	3	4	20,20,21	0.59	0	24,28,31	1.00	1 (4%)
3	NAG	O	1	3,1	14,14,15	0.51	0	17,19,21	0.84	0
3	NAG	O	2	3	14,14,15	0.54	0	17,19,21	1.01	2 (11%)
5	GAL	P	1	5	11,11,12	0.67	0	15,15,17	1.04	0
5	SIA	P	2	5	20,20,21	0.63	0	24,28,31	0.89	1 (4%)
5	GAL	Q	1	5	11,11,12	0.61	0	15,15,17	0.67	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\|>2$	Counts	RMSZ	$\#\|Z\|>2$
5	SIA	Q	2	5	20,20,21	0.61	0	24,28,31	0.91	1 (4%)
3	NAG	R	1	3,1	14,14,15	0.52	0	17,19,21	0.86	0
3	NAG	R	2	3	14,14,15	0.55	0	17,19,21	0.68	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	M	1	3,1	-	0/6/23/26	0/1/1/1
3	NAG	M	2	3	-	1/6/23/26	0/1/1/1
4	NAG	N	1	4	-	2/6/26/26	0/1/1/1
4	GAL	N	2	4	-	2/2/19/22	0/1/1/1
4	SIA	N	3	4	-	0/18/34/38	0/1/1/1
3	NAG	O	1	3,1	-	0/6/23/26	0/1/1/1
3	NAG	O	2	3	-	2/6/23/26	0/1/1/1
5	GAL	P	1	5	-	2/2/19/22	0/1/1/1
5	SIA	P	2	5	-	0/18/34/38	0/1/1/1
5	GAL	Q	1	5	-	1/2/19/22	0/1/1/1
5	SIA	Q	2	5	-	0/18/34/38	0/1/1/1
3	NAG	R	1	3,1	-	0/6/23/26	0/1/1/1
3	NAG	R	2	3	-	2/6/23/26	0/1/1/1

There are no bond length outliers.

The worst 5 of 10 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
4	N	2	GAL	O3-C3-C2	-3.41	103.47	109.99
5	P	2	SIA	C8-C7-C6	-2.37	108.55	113.03
3	M	1	NAG	O5-C1-C2	-2.33	107.60	111.29
3	O	2	NAG	O5-C1-C2	-2.27	107.70	111.29
5	Q	2	SIA	C6-C5-N5	-2.23	107.22	110.91

There are no chirality outliers.

5 of 12 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	O	2	NAG	C8-C7-N2-C2
3	O	2	NAG	O7-C7-N2-C2

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms
5	P	1	GAL	O5-C5-C6-O6
4	N	1	NAG	O5-C5-C6-O6
3	R	2	NAG	C8-C7-N2-C2

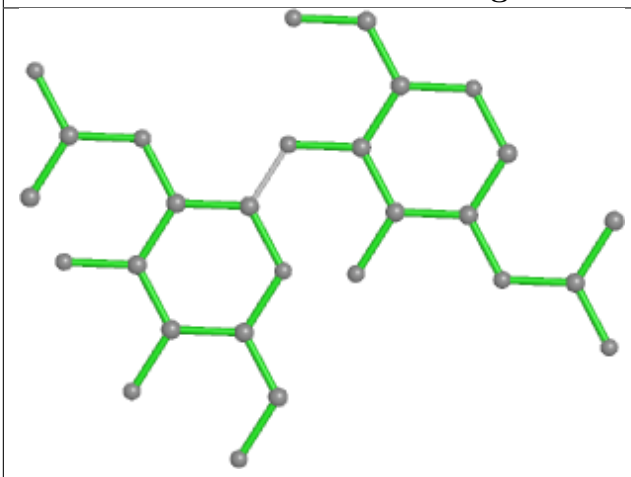
There are no ring outliers.

5 monomers are involved in 6 short contacts:

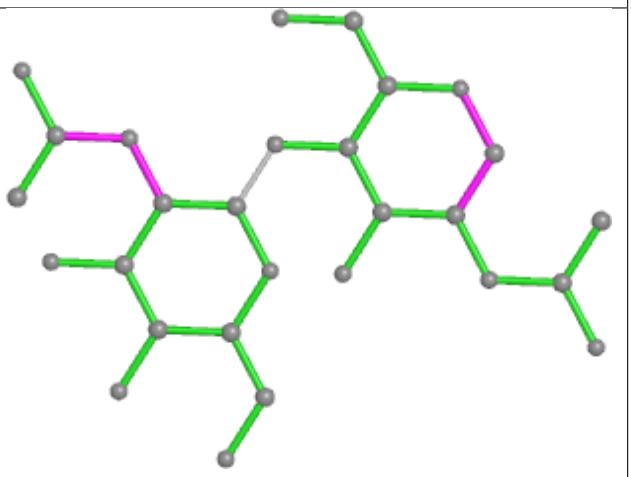
Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	P	2	SIA	2	0
3	R	1	NAG	1	0
5	Q	2	SIA	1	0
3	R	2	NAG	1	0
4	N	3	SIA	1	0

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

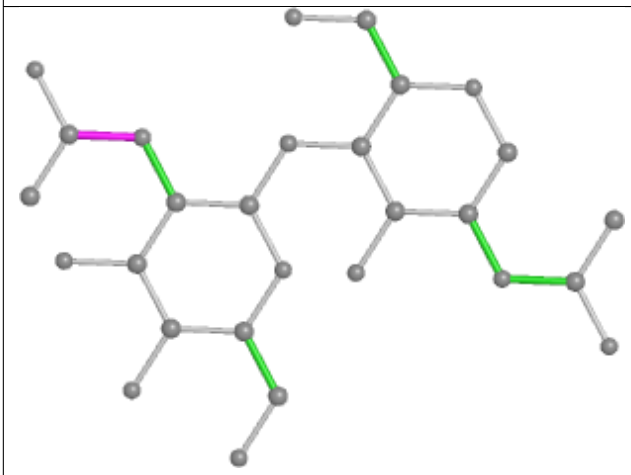
Oligosaccharide Chain M



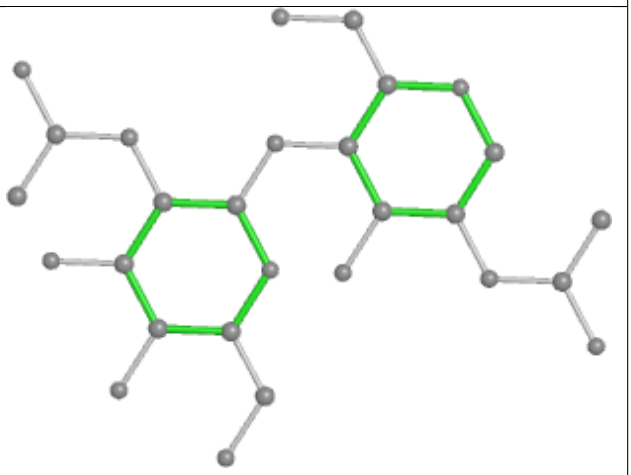
Bond lengths



Bond angles

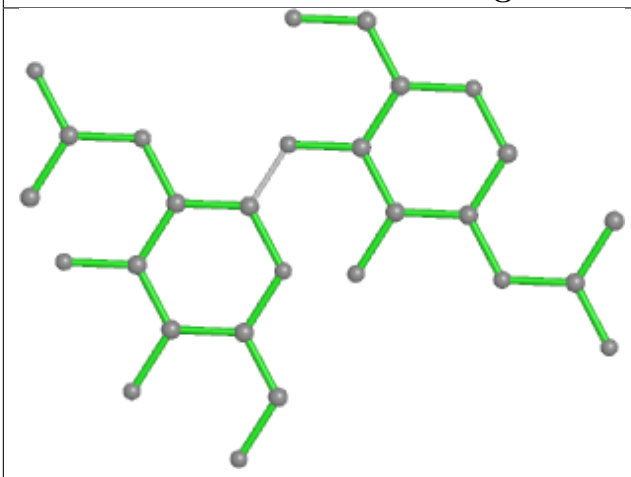


Torsions

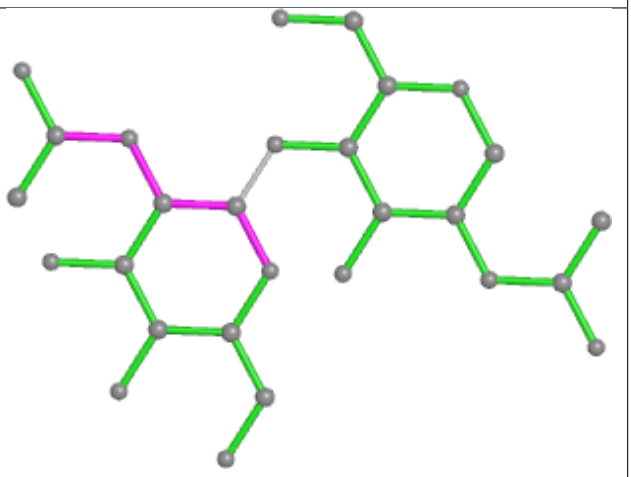


Rings

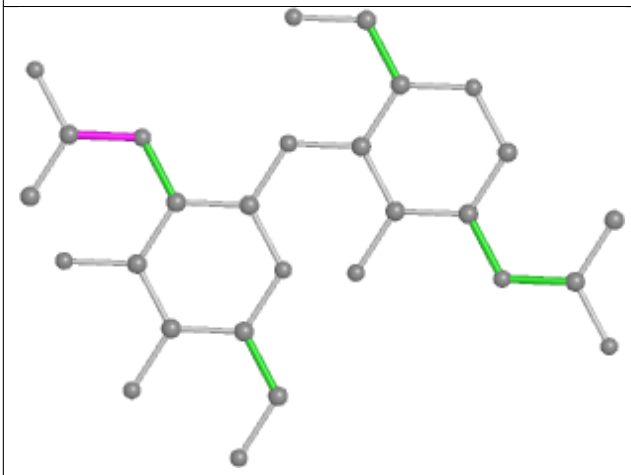
Oligosaccharide Chain O



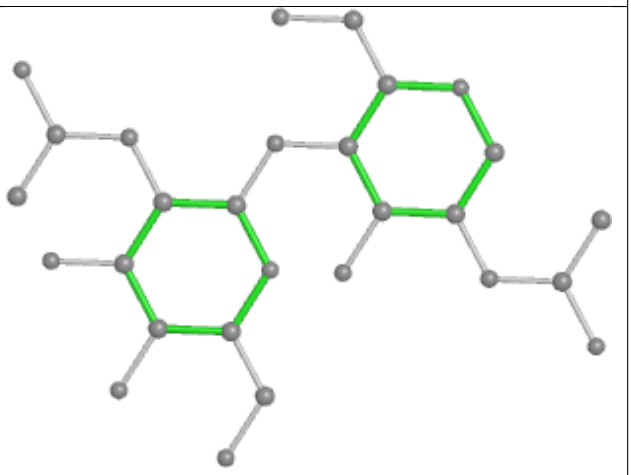
Bond lengths



Bond angles

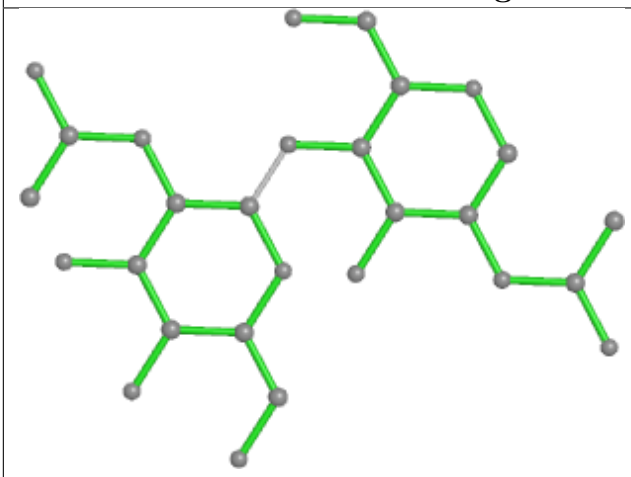


Torsions

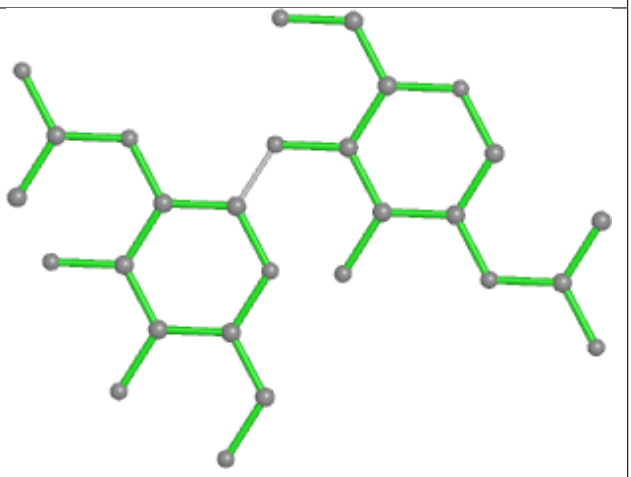


Rings

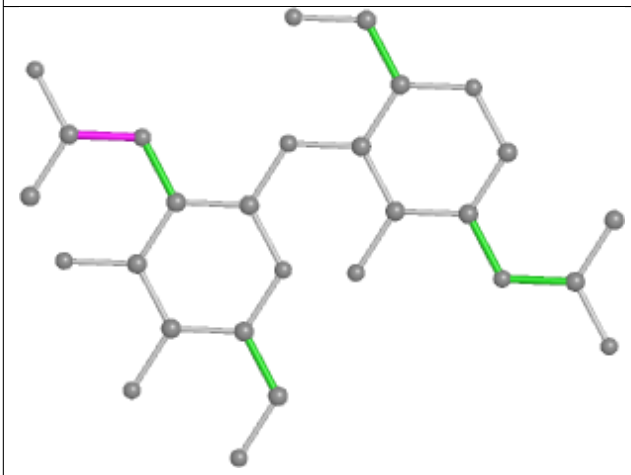
Oligosaccharide Chain R



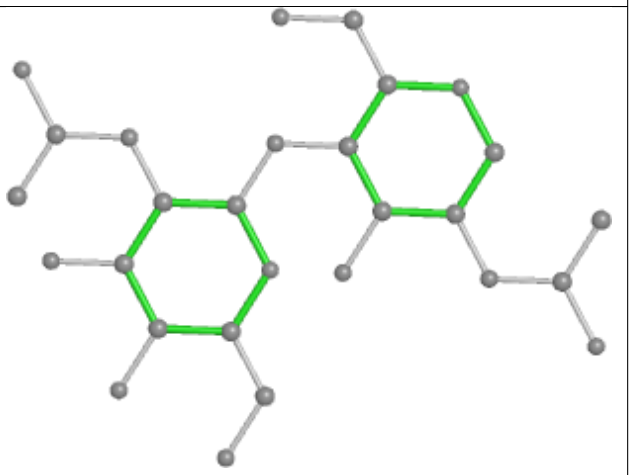
Bond lengths



Bond angles

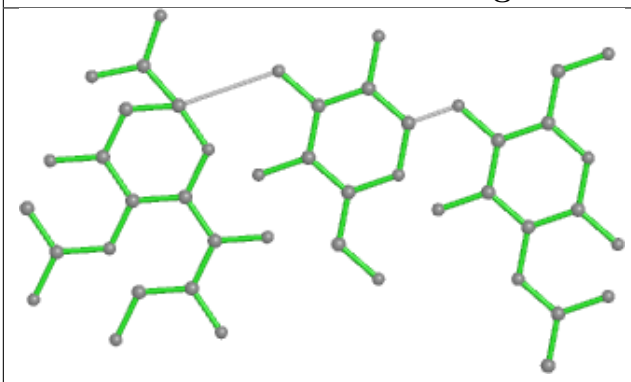


Torsions

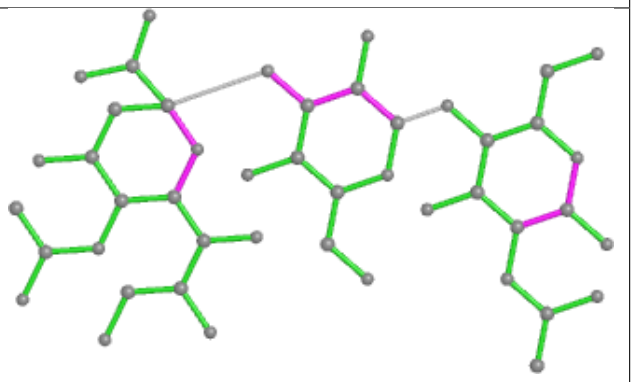


Rings

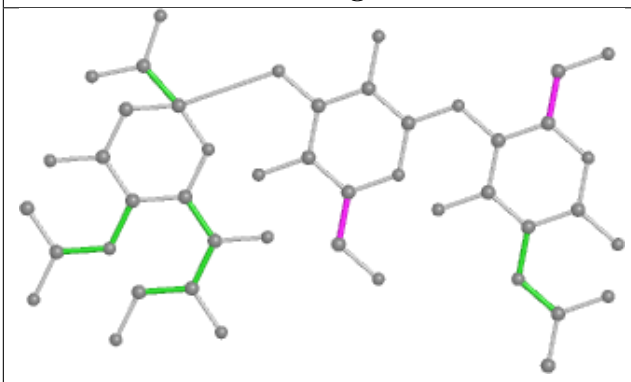
Oligosaccharide Chain N



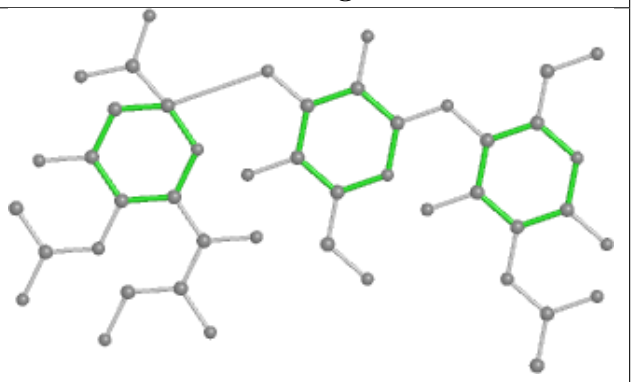
Bond lengths



Bond angles

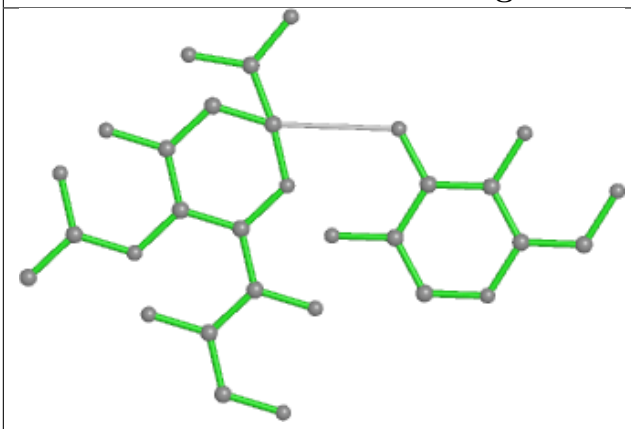


Torsions

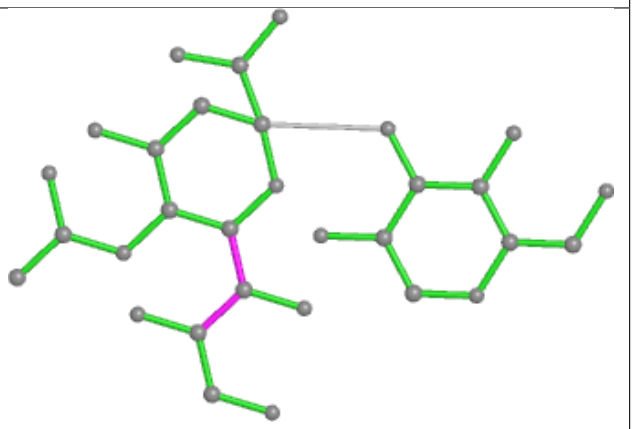


Rings

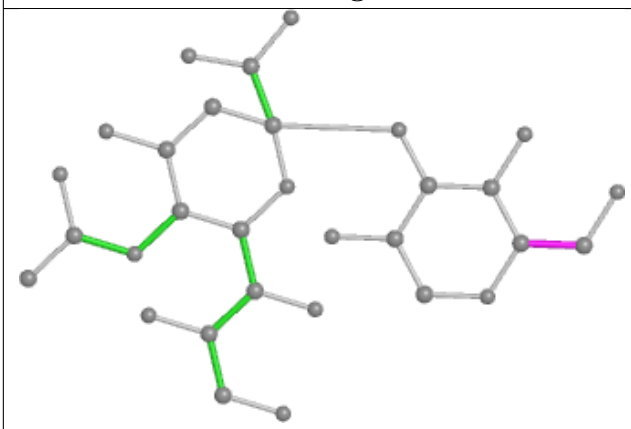
Oligosaccharide Chain P



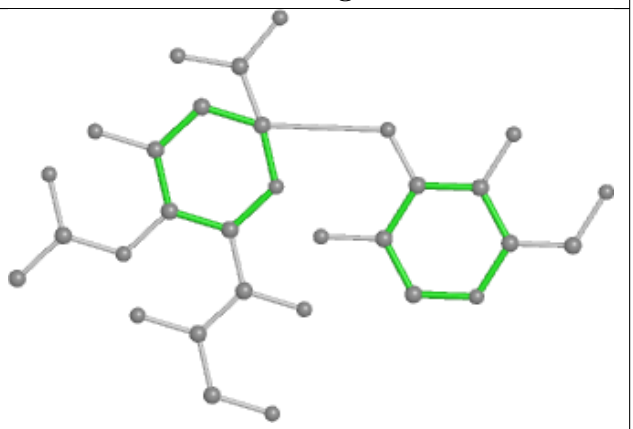
Bond lengths



Bond angles

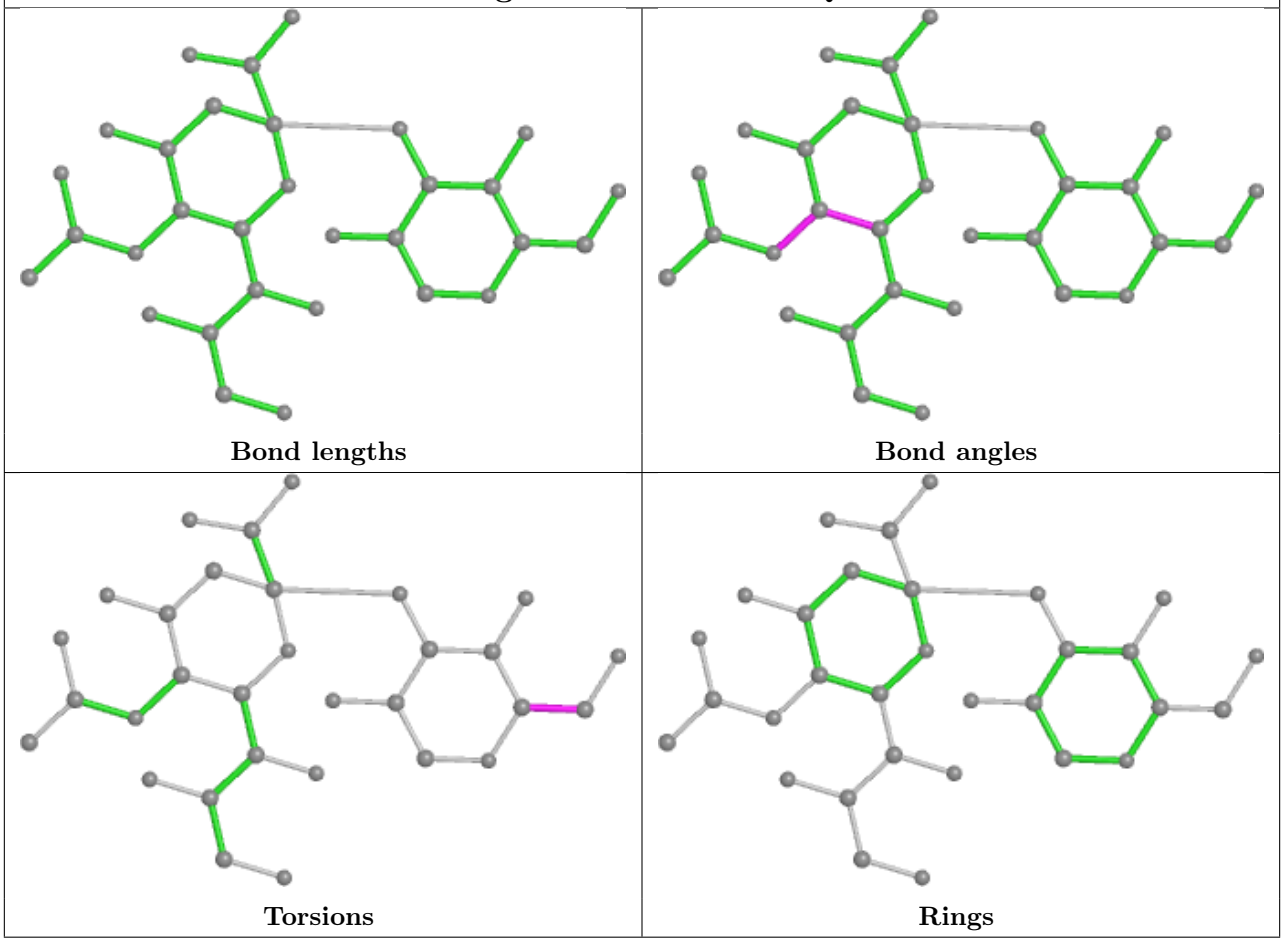


Torsions



Rings

Oligosaccharide Chain Q



5.6 Ligand geometry (i)

12 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	K	641	1	14,14,15	0.51	0	17,19,21	0.74	0
6	NAG	I	531	1	14,14,15	0.48	0	17,19,21	1.06	1 (5%)
6	NAG	I	541	1	14,14,15	0.47	0	17,19,21	0.92	0
6	NAG	G	451	1	14,14,15	0.50	0	17,19,21	0.75	1 (5%)
6	NAG	I	521	1	14,14,15	0.47	0	17,19,21	0.82	1 (5%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\|>2$	Counts	RMSZ	$\#\|Z\|>2$
6	NAG	K	651	1	14,14,15	0.46	0	17,19,21	0.79	0
6	NAG	E	341	1	14,14,15	0.52	0	17,19,21	1.01	1 (5%)
6	NAG	K	621	1	14,14,15	0.44	0	17,19,21	0.91	2 (11%)
6	NAG	G	411	1	14,14,15	0.49	0	17,19,21	0.90	1 (5%)
6	NAG	D	261	2	14,14,15	0.47	0	17,19,21	0.91	1 (5%)
6	NAG	G	431	1	14,14,15	0.53	0	17,19,21	0.97	2 (11%)
6	NAG	A	332	1	14,14,15	0.55	0	17,19,21	0.74	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	K	641	1	-	0/6/23/26	0/1/1/1
6	NAG	I	531	1	-	1/6/23/26	0/1/1/1
6	NAG	I	541	1	-	0/6/23/26	0/1/1/1
6	NAG	G	451	1	-	0/6/23/26	0/1/1/1
6	NAG	I	521	1	-	2/6/23/26	0/1/1/1
6	NAG	K	651	1	-	0/6/23/26	0/1/1/1
6	NAG	E	341	1	-	1/6/23/26	0/1/1/1
6	NAG	K	621	1	-	0/6/23/26	0/1/1/1
6	NAG	G	411	1	-	2/6/23/26	0/1/1/1
6	NAG	D	261	2	-	0/6/23/26	0/1/1/1
6	NAG	G	431	1	-	0/6/23/26	0/1/1/1
6	NAG	A	332	1	-	2/6/23/26	0/1/1/1

There are no bond length outliers.

The worst 5 of 10 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
6	I	531	NAG	C1-O5-C5	2.93	116.16	112.19
6	G	411	NAG	C1-O5-C5	2.41	115.46	112.19
6	I	521	NAG	O5-C5-C6	2.38	110.93	107.20
6	G	431	NAG	C1-O5-C5	2.29	115.30	112.19
6	G	431	NAG	C2-N2-C7	-2.22	119.74	122.90

There are no chirality outliers.

5 of 8 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
6	G	411	NAG	O5-C5-C6-O6
6	I	521	NAG	C8-C7-N2-C2
6	I	521	NAG	O7-C7-N2-C2
6	G	411	NAG	C4-C5-C6-O6
6	A	332	NAG	C8-C7-N2-C2

There are no ring outliers.

3 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	G	451	NAG	2	0
6	G	431	NAG	1	0
6	A	332	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	323/329 (98%)	-0.10	8 (2%) 57 56	22, 35, 56, 75	0
1	C	321/329 (97%)	-0.16	5 (1%) 72 70	27, 42, 58, 76	0
1	E	321/329 (97%)	-0.15	4 (1%) 79 78	24, 35, 51, 69	0
1	G	323/329 (98%)	-0.16	2 (0%) 89 88	24, 36, 50, 70	0
1	I	319/329 (96%)	-0.28	4 (1%) 77 76	24, 36, 54, 76	0
1	K	318/329 (96%)	-0.11	2 (0%) 89 88	27, 40, 58, 74	0
2	B	169/177 (95%)	0.88	27 (15%) 1 1	23, 45, 90, 106	0
2	D	170/177 (96%)	0.17	6 (3%) 44 43	25, 46, 63, 75	0
2	F	170/177 (96%)	0.04	3 (1%) 68 66	24, 40, 60, 67	0
2	H	172/177 (97%)	-0.04	5 (2%) 51 50	24, 43, 58, 68	0
2	J	170/177 (96%)	0.70	34 (20%) 1 0	24, 46, 78, 83	0
2	L	169/177 (95%)	0.47	22 (13%) 3 3	26, 45, 81, 94	0
All	All	2945/3036 (97%)	0.02	122 (4%) 37 36	22, 39, 68, 106	0

The worst 5 of 122 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	L	168	LEU	8.7
2	B	166	ALA	7.7
2	B	168	LEU	7.2
2	B	160	PRO	7.0
2	J	147	THR	6.6

6.2 Non-standard residues in protein, DNA, RNA chains

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

6.3 Carbohydrates (i)

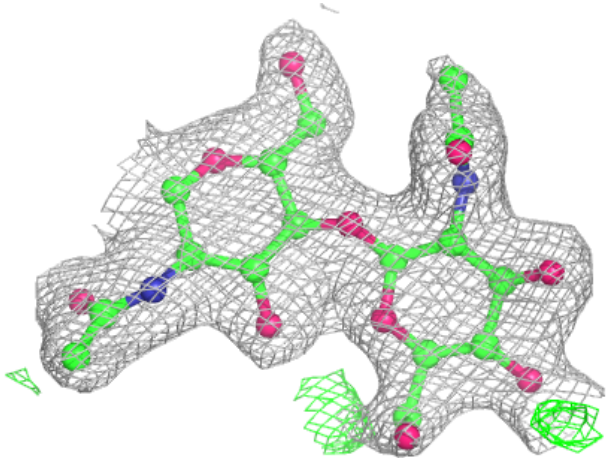
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95^{th} percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9'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

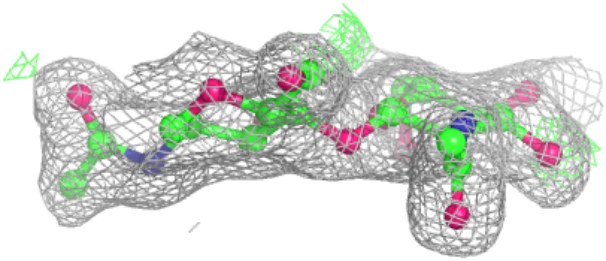
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
3	NAG	O	2	14/15	0.77	0.35	62,77,84,85	0
3	NAG	O	1	14/15	0.79	0.25	55,63,71,71	0
3	NAG	R	2	14/15	0.79	0.19	67,75,84,85	0
4	GAL	N	2	11/12	0.82	0.12	53,65,70,76	0
4	NAG	N	1	15/15	0.83	0.27	74,83,91,94	0
5	GAL	P	1	11/12	0.86	0.12	55,60,68,69	0
5	GAL	Q	1	11/12	0.86	0.19	59,70,77,77	0
3	NAG	M	2	14/15	0.89	0.15	44,53,62,67	0
5	SIA	Q	2	20/21	0.89	0.12	42,51,57,57	0
5	SIA	P	2	20/21	0.91	0.11	36,45,49,51	0
3	NAG	R	1	14/15	0.92	0.12	51,58,66,69	0
4	SIA	N	3	20/21	0.94	0.10	33,43,47,50	0
3	NAG	M	1	14/15	0.96	0.08	32,39,46,48	0

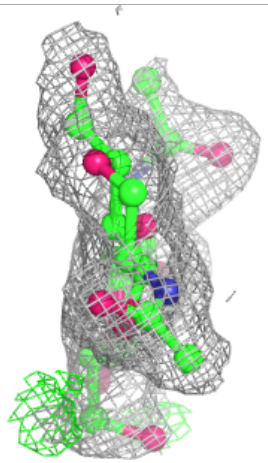
The following is a graphical depiction of the model fit to experimental electron density for oligosaccharide. Each fit is shown from different orientation to approximate a three-dimensional view.

Electron density around Chain M:

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

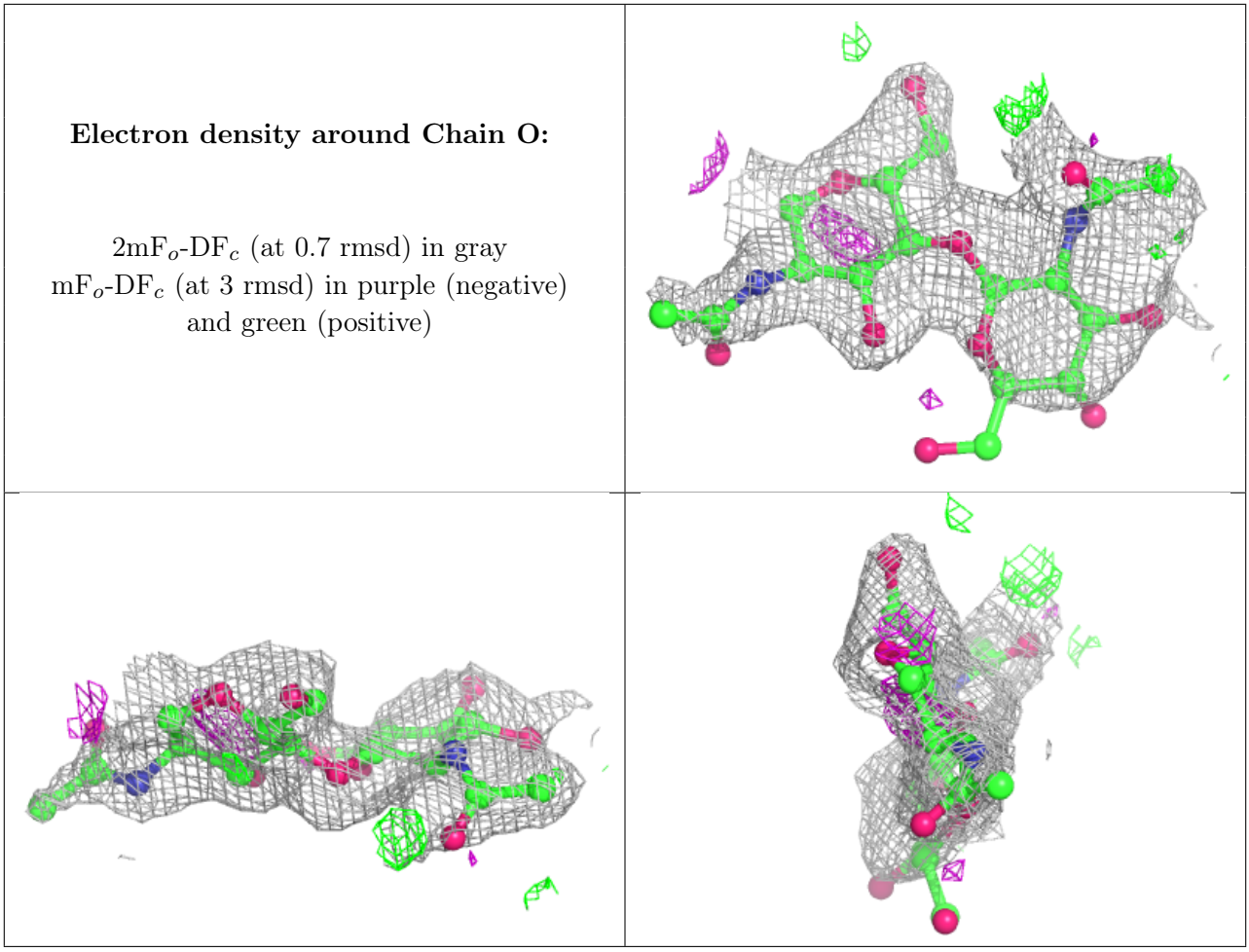






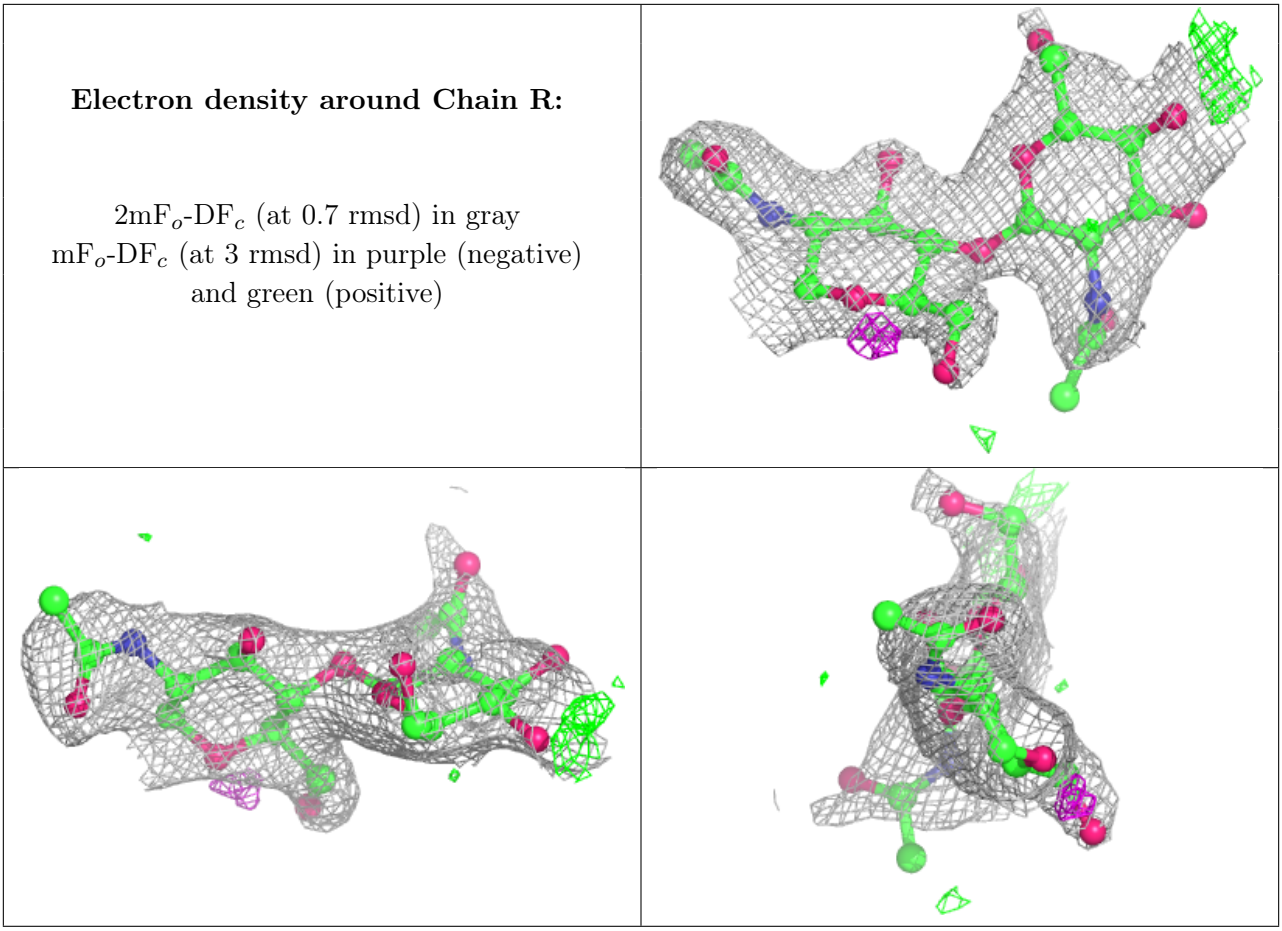
Electron density around Chain O:

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



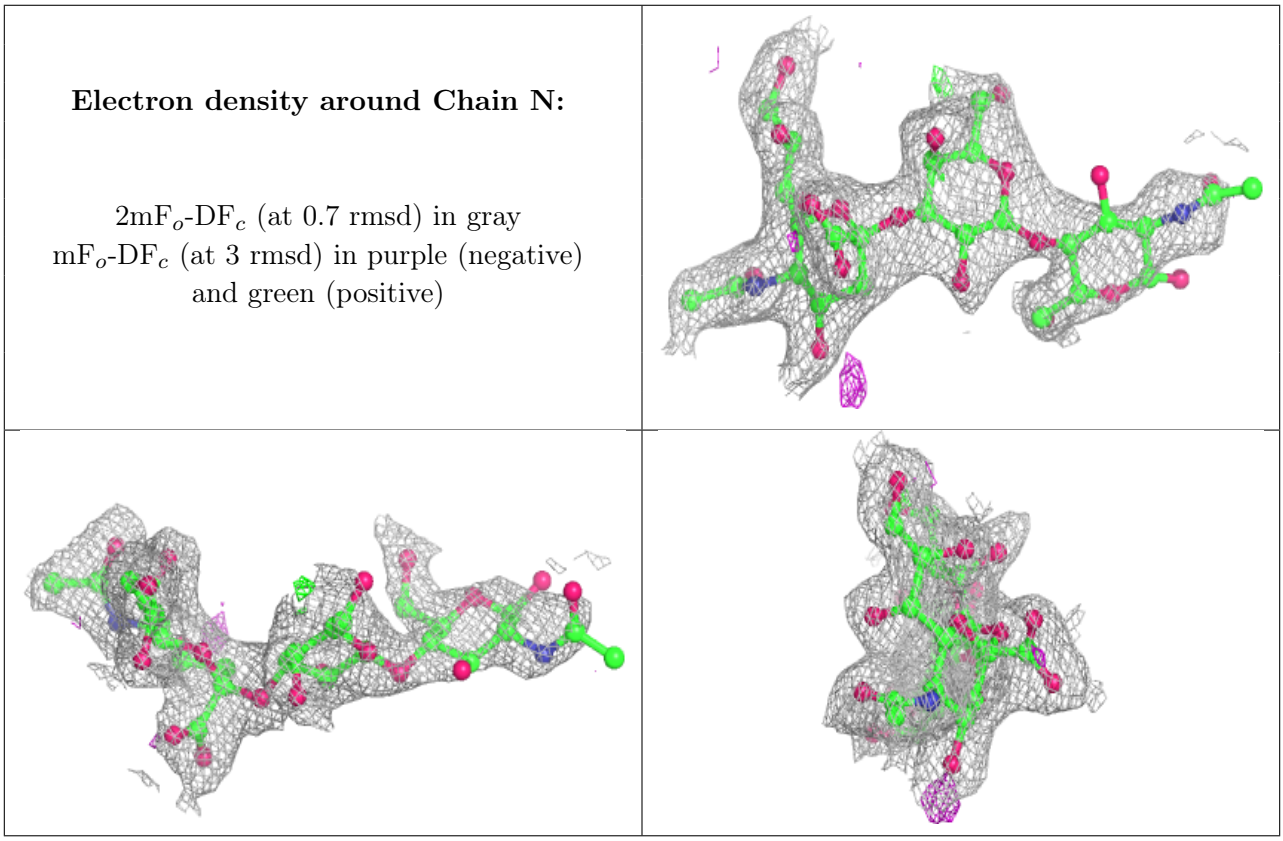
Electron density around Chain R:

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



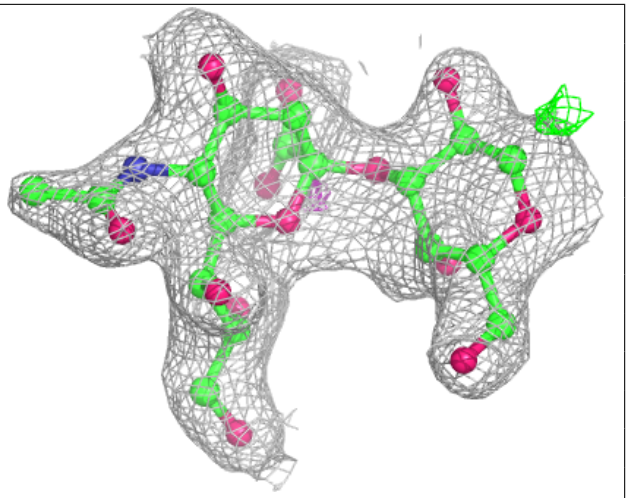
Electron density around Chain N:

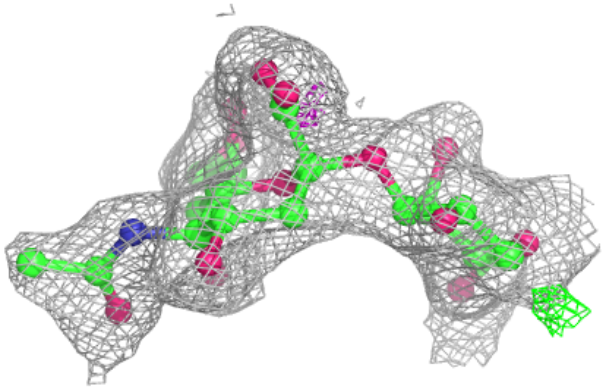
$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

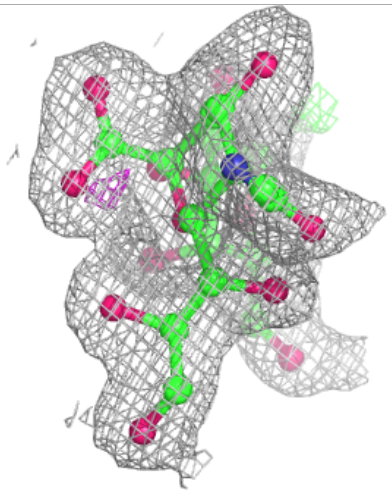


Electron density around Chain P:

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

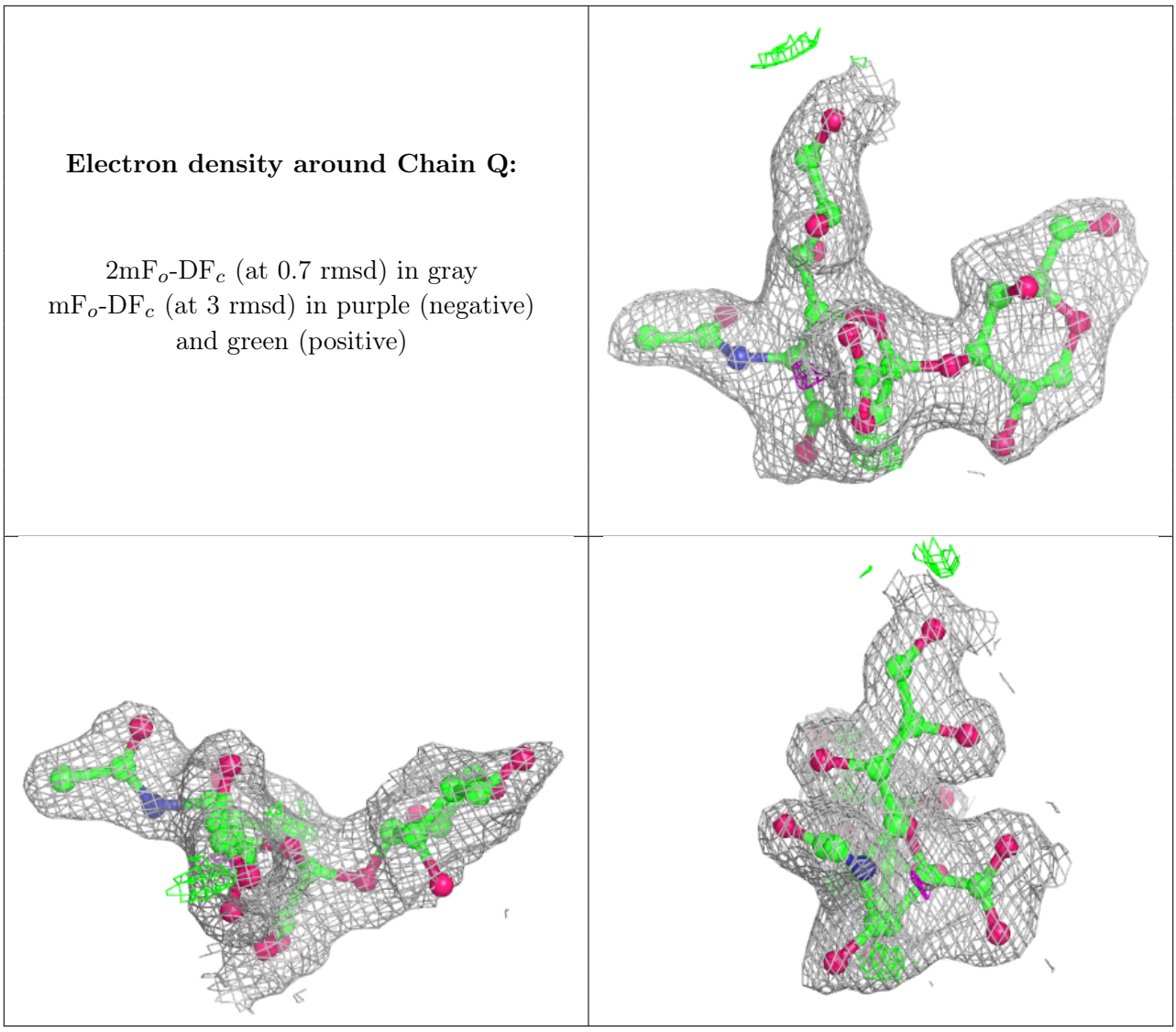






Electron density around Chain Q:

$2mF_o$-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
6	NAG	I	521	14/15	0.68	0.20	73,83,88,92	0
6	NAG	G	451	14/15	0.69	0.37	65,84,90,94	0
6	NAG	E	341	14/15	0.74	0.32	76,79,87,89	0
6	NAG	K	651	14/15	0.75	0.22	53,74,82,85	0
6	NAG	A	332	14/15	0.77	0.37	51,62,69,70	0
6	NAG	K	621	14/15	0.80	0.18	53,67,77,83	0
6	NAG	K	641	14/15	0.81	0.21	57,63,74,78	0

Continued on next page...

Continued from previous page...

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
6	NAG	I	531	14/15	0.82	0.15	41,53,63,64	0
6	NAG	I	541	14/15	0.85	0.15	57,67,75,80	0
6	NAG	G	411	14/15	0.85	0.31	70,75,82,84	0
6	NAG	D	261	14/15	0.86	0.13	65,68,72,76	0
6	NAG	G	431	14/15	0.92	0.10	40,52,62,64	0

6.5 Other polymers (i)

There are no such residues in this entry.