



wwPDB X-ray Structure Validation Summary Report ⓘ

Aug 8, 2020 – 05:42 PM BST

PDB ID : 6N4D
Title : The crystal structure of neuramindase from A/canine/IL/11613/2015 (H3N2) influenza virus.
Authors : Yang, H.; Stevens, J.
Deposited on : 2018-11-19
Resolution : 1.80 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.13.1
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) : 2.13.1

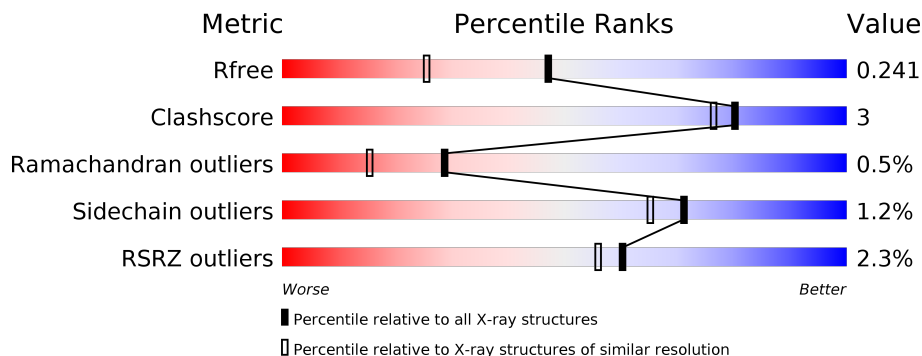
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.80 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	5950 (1.80-1.80)
Clashscore	141614	6793 (1.80-1.80)
Ramachandran outliers	138981	6697 (1.80-1.80)
Sidechain outliers	138945	6696 (1.80-1.80)
RSRZ outliers	127900	5850 (1.80-1.80)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	397	 2% 90% 8% ..
1	B	397	 4% 92% 6% ..
1	C	397	 2% 92% 6% ..
1	D	397	 % 90% 7% ..
2	E	2	 100%
2	G	2	 100%

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Mol	Chain	Length	Quality of chain
2	I	2	 100%
2	K	2	 100%
3	F	6	 100%
3	H	6	 67% 33%
3	J	6	 67% 33%
3	L	6	 100%

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	NAG	K	1	X	-	-	-

2 Entry composition i

There are 6 unique types of molecules in this entry. The entry contains 12810 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 Neuraminidase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	388	3017	1862	545	586	24	0	1	0
1	B	388	3017	1862	545	586	24	0	1	0
1	C	388	3017	1862	545	586	24	0	1	0
1	D	388	3017	1862	545	586	24	0	1	0

There are 36 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	73	GLY	-	expression tag	UNP A0A0H3YBU9
A	74	SER	-	expression tag	UNP A0A0H3YBU9
A	75	GLY	-	expression tag	UNP A0A0H3YBU9
A	76	ASP	-	expression tag	UNP A0A0H3YBU9
A	77	SER	-	expression tag	UNP A0A0H3YBU9
A	78	GLY	-	expression tag	UNP A0A0H3YBU9
A	79	SER	-	expression tag	UNP A0A0H3YBU9
A	80	PRO	-	expression tag	UNP A0A0H3YBU9
A	81	GLY	-	expression tag	UNP A0A0H3YBU9
B	73	GLY	-	expression tag	UNP A0A0H3YBU9
B	74	SER	-	expression tag	UNP A0A0H3YBU9
B	75	GLY	-	expression tag	UNP A0A0H3YBU9
B	76	ASP	-	expression tag	UNP A0A0H3YBU9
B	77	SER	-	expression tag	UNP A0A0H3YBU9
B	78	GLY	-	expression tag	UNP A0A0H3YBU9
B	79	SER	-	expression tag	UNP A0A0H3YBU9
B	80	PRO	-	expression tag	UNP A0A0H3YBU9
B	81	GLY	-	expression tag	UNP A0A0H3YBU9
C	73	GLY	-	expression tag	UNP A0A0H3YBU9
C	74	SER	-	expression tag	UNP A0A0H3YBU9
C	75	GLY	-	expression tag	UNP A0A0H3YBU9

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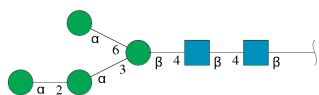
Chain	Residue	Modelled	Actual	Comment	Reference
C	76	ASP	-	expression tag	UNP A0A0H3YBU9
C	77	SER	-	expression tag	UNP A0A0H3YBU9
C	78	GLY	-	expression tag	UNP A0A0H3YBU9
C	79	SER	-	expression tag	UNP A0A0H3YBU9
C	80	PRO	-	expression tag	UNP A0A0H3YBU9
C	81	GLY	-	expression tag	UNP A0A0H3YBU9
D	73	GLY	-	expression tag	UNP A0A0H3YBU9
D	74	SER	-	expression tag	UNP A0A0H3YBU9
D	75	GLY	-	expression tag	UNP A0A0H3YBU9
D	76	ASP	-	expression tag	UNP A0A0H3YBU9
D	77	SER	-	expression tag	UNP A0A0H3YBU9
D	78	GLY	-	expression tag	UNP A0A0H3YBU9
D	79	SER	-	expression tag	UNP A0A0H3YBU9
D	80	PRO	-	expression tag	UNP A0A0H3YBU9
D	81	GLY	-	expression tag	UNP A0A0H3YBU9

- Molecule 2 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
			Total	C	N	O			
2	E	2	28	16	2	10	0	0	0
2	G	2	28	16	2	10	0	0	0
2	I	2	28	16	2	10	0	0	0
2	K	2	28	16	2	10	0	0	0

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

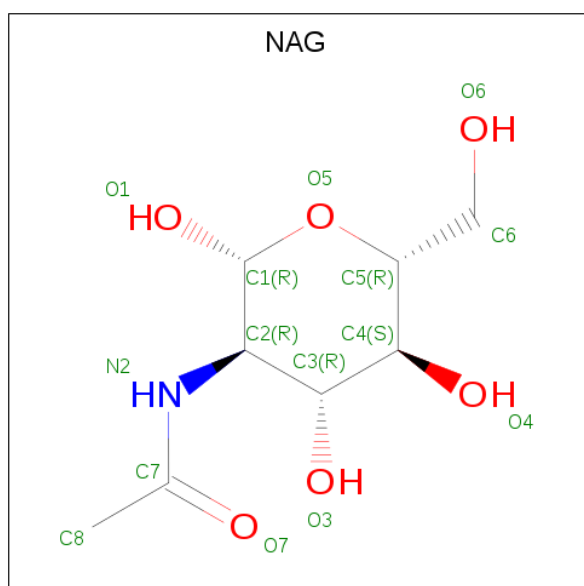


Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
3	F	6	Total	C	N	O	0	0	0
			72	40	2	30			
3	H	6	Total	C	N	O	0	0	0
			72	40	2	30			
3	J	6	Total	C	N	O	0	0	0
			72	40	2	30			
3	L	6	Total	C	N	O	0	0	0
			72	40	2	30			

- Molecule 4 is CALCIUM ION (three-letter code: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	B	1	Total	Ca	0	0
			1	1		
4	A	1	Total	Ca	0	0
			1	1		
4	D	1	Total	Ca	0	0
			1	1		
4	C	1	Total	Ca	0	0
			1	1		

- Molecule 5 is 2-acetamido-2-deoxy-beta-D-glucopyranose (three-letter code: NAG) (formula: C₈H₁₅NO₆).



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

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

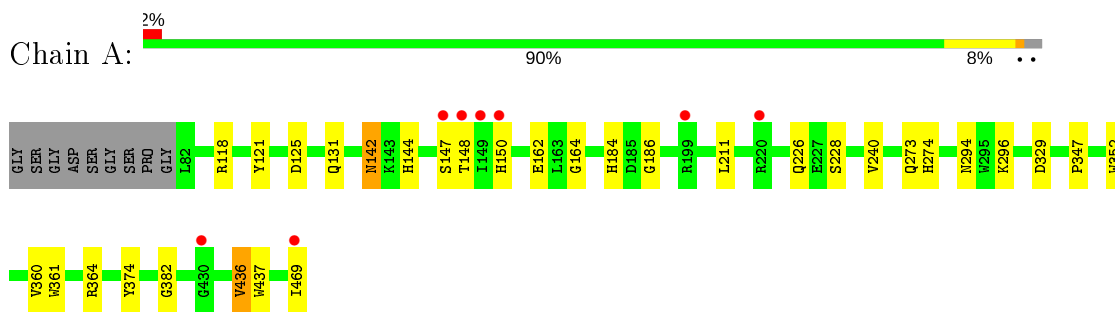
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	97	Total	O	0	0
			97	97		
6	B	40	Total	O	0	0
			40	40		
6	C	48	Total	O	0	0
			48	48		
6	D	97	Total	O	0	0
			97	97		

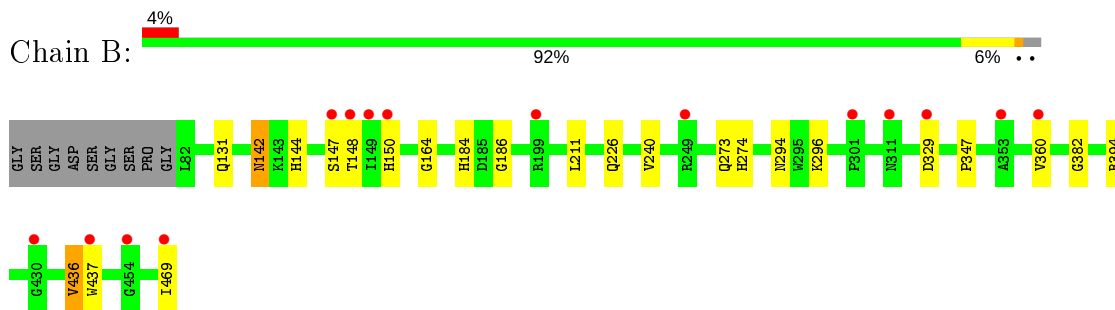
3 Residue-property plots [i](#)

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.

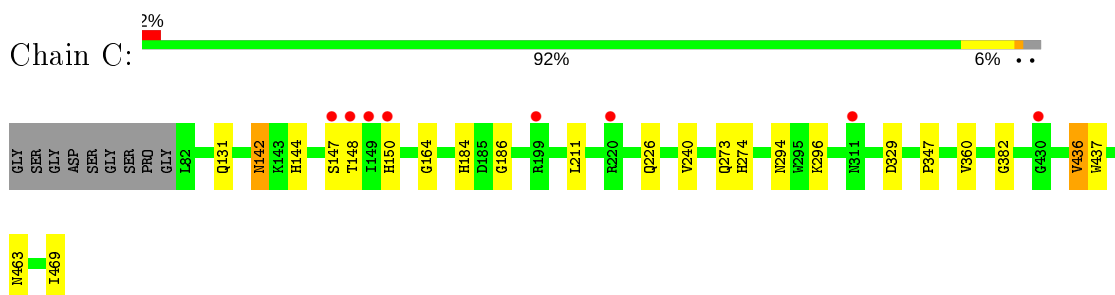
- Molecule 1: Neuraminidase



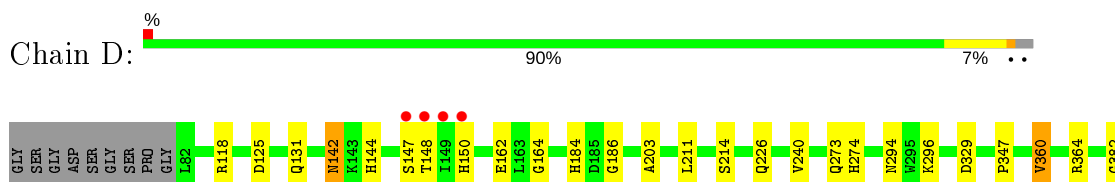
- Molecule 1: Neuraminidase

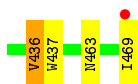


- Molecule 1: Neuraminidase



- Molecule 1: Neuraminidase





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

Chain E: 100%



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

Chain G: 100%



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

Chain I: 100%



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

Chain K: 100%



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

Chain F: 100%



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

Chain H: 67% 33%

MAG1
MAG2
BMA3
MAN4
MAN5
MAN6

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

Chain J:  67% 33%

MAG1
MAG2
BMA3
MAN4
MAN5
MAN6

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

Chain L:  100%

MAG1
MAG2
BMA3
MAN4
MAN5
MAN6

4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α , β , γ	110.45Å 110.56Å 126.62Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	49.43 – 1.80 35.39 – 1.80	Depositor EDS
% Data completeness (in resolution range)	99.7 (49.43-1.80) 99.4 (35.39-1.80)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.51 (at 1.81Å)	Xtrriage
Refinement program	REFMAC 5.8.0049	Depositor
R, R_{free}	0.210 , 0.238 0.216 , 0.241	Depositor DCC
R_{free} test set	7193 reflections (5.02%)	wwPDB-VP
Wilson B-factor (Å ²)	23.9	Xtrriage
Anisotropy	0.170	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 16.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.29$	Xtrriage
Estimated twinning fraction	0.467 for k,h,-l	Xtrriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	12810	wwPDB-VP
Average B, all atoms (Å ²)	28.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 35.03 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 6.2179e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

¹ Intensities estimated from amplitudes.

² Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: CA, BMA, NAG, MAN

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	0.99	2/3089 (0.1%)	0.92	4/4195 (0.1%)
1	B	0.82	0/3089	0.86	2/4195 (0.0%)
1	C	0.80	0/3089	0.86	1/4195 (0.0%)
1	D	0.99	1/3089 (0.0%)	0.91	5/4195 (0.1%)
All	All	0.90	3/12356 (0.0%)	0.89	12/16780 (0.1%)

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	162	GLU	CD-OE1	-5.95	1.19	1.25
1	A	162	GLU	CD-OE1	-5.12	1.20	1.25
1	A	361	TRP	CE3-CZ3	5.01	1.47	1.38

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

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	436	VAL	N-CA-C	6.46	128.44	111.00
1	D	436	VAL	N-CA-C	6.46	128.44	111.00
1	B	436	VAL	N-CA-C	6.21	127.76	111.00
1	C	436	VAL	N-CA-C	6.09	127.46	111.00
1	D	125	ASP	CB-CG-OD1	5.78	123.50	118.30

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	3017	0	2837	18	0
1	B	3017	0	2837	15	0
1	C	3017	0	2837	15	0
1	D	3017	0	2837	17	0
2	E	28	0	25	1	0
2	G	28	0	25	0	0
2	I	28	0	25	0	0
2	K	28	0	25	0	0
3	F	72	0	61	0	0
3	H	72	0	61	1	0
3	J	72	0	61	1	0
3	L	72	0	61	0	0
4	A	1	0	0	0	0
4	B	1	0	0	0	0
4	C	1	0	0	0	0
4	D	1	0	0	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	0	0
6	A	97	0	0	0	0
6	B	40	0	0	0	0
6	C	48	0	0	0	0
6	D	97	0	0	0	0
All	All	12810	0	11744	66	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 3.

The worst 5 of 66 close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:226:GLN:HE21	1:D:240:VAL:H	1.34	0.75
1:A:226:GLN:HE21	1:A:240:VAL:H	1.35	0.73
1:A:144:HIS:HE2	1:D:463:ASN:H	1.38	0.72
1:B:144:HIS:HE2	1:C:463:ASN:H	1.40	0.70
1:C:226:GLN:HE21	1:C:240:VAL:H	1.40	0.69

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	387/397 (98%)	366 (95%)	19 (5%)	2 (0%)	29	15
1	B	387/397 (98%)	366 (95%)	19 (5%)	2 (0%)	29	15
1	C	387/397 (98%)	366 (95%)	19 (5%)	2 (0%)	29	15
1	D	387/397 (98%)	366 (95%)	19 (5%)	2 (0%)	29	15
All	All	1548/1588 (98%)	1464 (95%)	76 (5%)	8 (0%)	29	15

5 of 8 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	347	PRO
1	B	347	PRO
1	C	347	PRO
1	D	347	PRO
1	A	437	TRP

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	339/343 (99%)	335 (99%)	4 (1%)	71	65
1	B	339/343 (99%)	335 (99%)	4 (1%)	71	65
1	C	339/343 (99%)	335 (99%)	4 (1%)	71	65
1	D	339/343 (99%)	335 (99%)	4 (1%)	71	65
All	All	1356/1372 (99%)	1340 (99%)	16 (1%)	71	65

5 of 16 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	329	ASP
1	C	142	ASN
1	D	142	ASN
1	B	211	LEU
1	D	148	THR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 57 such sidechains are listed below:

Mol	Chain	Res	Type
1	B	393	ASN
1	C	142	ASN
1	D	356	ASN
1	B	419	ASN
1	C	104	ASN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

32 non-standard protein/DNA/RNA residues 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
2	NAG	K	2	2	14,14,15	0.59	0	17,19,21	1.31	1 (5%)
3	MAN	J	6	3	11,11,12	1.22	1 (9%)	15,15,17	1.98	4 (26%)
3	MAN	L	6	3	11,11,12	0.84	0	15,15,17	1.93	3 (20%)
5	NAG	C	504	1	14,14,15	0.56	0	17,19,21	1.33	2 (11%)
3	MAN	L	5	3	11,11,12	1.20	1 (9%)	15,15,17	1.80	3 (20%)
3	NAG	L	1	1,3	14,14,15	1.08	2 (14%)	17,19,21	1.42	4 (23%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	MAN	L	4	3	11,11,12	1.54	3 (27%)	15,15,17	1.64	5 (33%)
2	NAG	E	1	1,2	14,14,15	0.71	0	17,19,21	2.30	7 (41%)
3	NAG	L	2	3	14,14,15	1.12	2 (14%)	17,19,21	1.82	4 (23%)
2	NAG	I	2	2	14,14,15	0.50	0	17,19,21	1.27	1 (5%)
2	NAG	G	2	2	14,14,15	0.52	0	17,19,21	1.04	1 (5%)
3	MAN	F	4	3	11,11,12	1.41	2 (18%)	15,15,17	1.37	2 (13%)
5	NAG	D	504	1	14,14,15	0.66	0	17,19,21	1.44	3 (17%)
3	NAG	F	2	3	14,14,15	1.22	2 (14%)	17,19,21	1.85	4 (23%)
3	NAG	H	1	1,3	14,14,15	0.74	0	17,19,21	1.53	3 (17%)
2	NAG	I	1	1,2	14,14,15	1.04	1 (7%)	17,19,21	0.81	0
3	MAN	F	6	3	11,11,12	0.90	0	15,15,17	2.13	4 (26%)
3	MAN	F	5	3	11,11,12	1.37	1 (9%)	15,15,17	1.88	3 (20%)
3	MAN	J	5	3	11,11,12	0.65	0	15,15,17	2.50	6 (40%)
3	NAG	H	2	3	14,14,15	0.94	1 (7%)	17,19,21	1.69	5 (29%)
3	MAN	J	4	3	11,11,12	0.85	0	15,15,17	1.64	2 (13%)
3	MAN	H	4	3	11,11,12	0.83	0	15,15,17	1.55	3 (20%)
5	NAG	B	504	1	14,14,15	0.51	0	17,19,21	1.25	2 (11%)
3	NAG	J	1	1,3	14,14,15	0.69	0	17,19,21	1.51	3 (17%)
3	NAG	F	1	1,3	14,14,15	1.09	1 (7%)	17,19,21	1.59	3 (17%)
3	MAN	H	6	3	11,11,12	0.97	0	15,15,17	1.66	3 (20%)
5	NAG	A	504	1	14,14,15	0.69	0	17,19,21	1.26	1 (5%)
2	NAG	E	2	2	14,14,15	0.60	0	17,19,21	1.28	2 (11%)
2	NAG	G	1	1,2	14,14,15	0.98	1 (7%)	17,19,21	0.76	0
3	MAN	H	5	3	11,11,12	0.75	0	15,15,17	2.28	7 (46%)
3	NAG	J	2	3	14,14,15	1.02	1 (7%)	17,19,21	2.04	7 (41%)
2	NAG	K	1	1,2	14,14,15	0.67	0	17,19,21	2.63	5 (29%)

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
2	NAG	K	2	2	-	1/6/23/26	0/1/1/1
3	MAN	J	6	3	-	2/2/19/22	0/1/1/1
3	MAN	L	6	3	-	2/2/19/22	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
5	NAG	C	504	1	-	2/6/23/26	0/1/1/1
3	MAN	L	5	3	-	0/2/19/22	0/1/1/1
3	NAG	L	1	1,3	-	0/6/23/26	0/1/1/1
3	MAN	L	4	3	-	0/2/19/22	0/1/1/1
2	NAG	E	1	1,2	-	0/6/23/26	0/1/1/1
3	NAG	L	2	3	-	0/6/23/26	0/1/1/1
2	NAG	I	2	2	-	0/6/23/26	0/1/1/1
2	NAG	G	2	2	-	0/6/23/26	0/1/1/1
3	MAN	F	4	3	-	0/2/19/22	0/1/1/1
5	NAG	D	504	1	-	0/6/23/26	0/1/1/1
3	NAG	F	2	3	-	0/6/23/26	0/1/1/1
3	NAG	H	1	1,3	-	0/6/23/26	0/1/1/1
2	NAG	I	1	1,2	-	0/6/23/26	0/1/1/1
3	MAN	F	6	3	-	2/2/19/22	0/1/1/1
3	MAN	F	5	3	-	0/2/19/22	0/1/1/1
3	MAN	J	5	3	-	1/2/19/22	0/1/1/1
3	NAG	H	2	3	-	0/6/23/26	0/1/1/1
3	MAN	J	4	3	-	0/2/19/22	0/1/1/1
3	MAN	H	4	3	-	0/2/19/22	0/1/1/1
5	NAG	B	504	1	-	2/6/23/26	0/1/1/1
3	NAG	J	1	1,3	-	0/6/23/26	0/1/1/1
3	NAG	F	1	1,3	-	0/6/23/26	0/1/1/1
3	MAN	H	6	3	-	2/2/19/22	0/1/1/1
5	NAG	A	504	1	-	0/6/23/26	0/1/1/1
2	NAG	E	2	2	-	0/6/23/26	0/1/1/1
2	NAG	G	1	1,2	-	0/6/23/26	0/1/1/1
3	MAN	H	5	3	-	2/2/19/22	0/1/1/1
3	NAG	J	2	3	-	0/6/23/26	0/1/1/1
2	NAG	K	1	1,2	1/1/5/7	0/6/23/26	0/1/1/1

The worst 5 of 19 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	L	4	MAN	O5-C1	-3.80	1.37	1.43
3	F	4	MAN	O5-C1	-3.65	1.37	1.43
3	F	5	MAN	O5-C1	-3.51	1.38	1.43
2	I	1	NAG	O5-C1	-3.44	1.38	1.43
3	F	2	NAG	O5-C1	-3.02	1.38	1.43

The worst 5 of 103 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	K	1	NAG	C1-O5-C5	6.54	121.05	112.19
3	J	5	MAN	C1-O5-C5	6.36	120.81	112.19
2	K	1	NAG	O5-C1-C2	6.32	121.27	111.29
3	J	6	MAN	O5-C5-C6	5.58	115.94	107.20
2	E	1	NAG	C1-O5-C5	5.16	119.19	112.19

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
2	K	1	NAG	C1

5 of 16 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	C	504	NAG	C4-C5-C6-O6
5	B	504	NAG	C4-C5-C6-O6
3	H	6	MAN	O5-C5-C6-O6
3	J	6	MAN	O5-C5-C6-O6
5	B	504	NAG	O5-C5-C6-O6

There are no ring outliers.

4 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	J	6	MAN	1	0
2	E	1	NAG	1	0
3	H	6	MAN	1	0
2	E	2	NAG	1	0

5.5 Carbohydrates [i](#)

32 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
2	NAG	E	1	1,2	14,14,15	0.71	0	17,19,21	2.30	7 (41%)
2	NAG	E	2	2	14,14,15	0.60	0	17,19,21	1.28	2 (11%)
3	NAG	F	1	1,3	14,14,15	1.09	1 (7%)	17,19,21	1.59	3 (17%)
3	NAG	F	2	3	14,14,15	1.22	2 (14%)	17,19,21	1.85	4 (23%)
3	BMA	F	3	3	11,11,12	0.96	1 (9%)	15,15,17	1.15	1 (6%)
3	MAN	F	4	3	11,11,12	1.41	2 (18%)	15,15,17	1.37	2 (13%)
3	MAN	F	5	3	11,11,12	1.37	1 (9%)	15,15,17	1.88	3 (20%)
3	MAN	F	6	3	11,11,12	0.90	0	15,15,17	2.13	4 (26%)
2	NAG	G	1	1,2	14,14,15	0.98	1 (7%)	17,19,21	0.76	0
2	NAG	G	2	2	14,14,15	0.52	0	17,19,21	1.04	1 (5%)
3	NAG	H	1	1,3	14,14,15	0.74	0	17,19,21	1.53	3 (17%)
3	NAG	H	2	3	14,14,15	0.94	1 (7%)	17,19,21	1.69	5 (29%)
3	BMA	H	3	3	11,11,12	0.68	0	15,15,17	1.54	2 (13%)
3	MAN	H	4	3	11,11,12	0.83	0	15,15,17	1.55	3 (20%)
3	MAN	H	5	3	11,11,12	0.75	0	15,15,17	2.28	7 (46%)
3	MAN	H	6	3	11,11,12	0.97	0	15,15,17	1.66	3 (20%)
2	NAG	I	1	1,2	14,14,15	1.04	1 (7%)	17,19,21	0.81	0
2	NAG	I	2	2	14,14,15	0.50	0	17,19,21	1.27	1 (5%)
3	NAG	J	1	1,3	14,14,15	0.69	0	17,19,21	1.51	3 (17%)
3	NAG	J	2	3	14,14,15	1.02	1 (7%)	17,19,21	2.04	7 (41%)
3	BMA	J	3	3	11,11,12	0.46	0	15,15,17	1.42	3 (20%)
3	MAN	J	4	3	11,11,12	0.85	0	15,15,17	1.64	2 (13%)
3	MAN	J	5	3	11,11,12	0.65	0	15,15,17	2.50	6 (40%)
3	MAN	J	6	3	11,11,12	1.22	1 (9%)	15,15,17	1.98	4 (26%)
2	NAG	K	1	1,2	14,14,15	0.67	0	17,19,21	2.63	5 (29%)
2	NAG	K	2	2	14,14,15	0.59	0	17,19,21	1.31	1 (5%)
3	NAG	L	1	1,3	14,14,15	1.08	2 (14%)	17,19,21	1.42	4 (23%)
3	NAG	L	2	3	14,14,15	1.12	2 (14%)	17,19,21	1.82	4 (23%)
3	BMA	L	3	3	11,11,12	1.24	2 (18%)	15,15,17	1.32	1 (6%)
3	MAN	L	4	3	11,11,12	1.54	3 (27%)	15,15,17	1.64	5 (33%)
3	MAN	L	5	3	11,11,12	1.20	1 (9%)	15,15,17	1.80	3 (20%)
3	MAN	L	6	3	11,11,12	0.84	0	15,15,17	1.93	3 (20%)

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
2	NAG	E	1	1,2	-	0/6/23/26	0/1/1/1
2	NAG	E	2	2	-	0/6/23/26	0/1/1/1
3	NAG	F	1	1,3	-	0/6/23/26	0/1/1/1
3	NAG	F	2	3	-	0/6/23/26	0/1/1/1
3	BMA	F	3	3	-	0/2/19/22	0/1/1/1
3	MAN	F	4	3	-	0/2/19/22	0/1/1/1
3	MAN	F	5	3	-	0/2/19/22	0/1/1/1
3	MAN	F	6	3	-	2/2/19/22	0/1/1/1
2	NAG	G	1	1,2	-	0/6/23/26	0/1/1/1
2	NAG	G	2	2	-	0/6/23/26	0/1/1/1
3	NAG	H	1	1,3	-	0/6/23/26	0/1/1/1
3	NAG	H	2	3	-	0/6/23/26	0/1/1/1
3	BMA	H	3	3	-	0/2/19/22	0/1/1/1
3	MAN	H	4	3	-	0/2/19/22	0/1/1/1
3	MAN	H	5	3	-	2/2/19/22	0/1/1/1
3	MAN	H	6	3	-	2/2/19/22	0/1/1/1
2	NAG	I	1	1,2	-	0/6/23/26	0/1/1/1
2	NAG	I	2	2	-	0/6/23/26	0/1/1/1
3	NAG	J	1	1,3	-	0/6/23/26	0/1/1/1
3	NAG	J	2	3	-	0/6/23/26	0/1/1/1
3	BMA	J	3	3	-	0/2/19/22	0/1/1/1
3	MAN	J	4	3	-	0/2/19/22	0/1/1/1
3	MAN	J	5	3	-	1/2/19/22	0/1/1/1
3	MAN	J	6	3	-	2/2/19/22	0/1/1/1
2	NAG	K	1	1,2	1/1/5/7	0/6/23/26	0/1/1/1
2	NAG	K	2	2	-	1/6/23/26	0/1/1/1
3	NAG	L	1	1,3	-	0/6/23/26	0/1/1/1
3	NAG	L	2	3	-	0/6/23/26	0/1/1/1
3	BMA	L	3	3	-	0/2/19/22	0/1/1/1
3	MAN	L	4	3	-	0/2/19/22	0/1/1/1
3	MAN	L	5	3	-	0/2/19/22	0/1/1/1
3	MAN	L	6	3	-	2/2/19/22	0/1/1/1

The worst 5 of 22 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	L	4	MAN	O5-C1	-3.80	1.37	1.43
3	F	4	MAN	O5-C1	-3.65	1.37	1.43
3	F	5	MAN	O5-C1	-3.51	1.38	1.43
2	I	1	NAG	O5-C1	-3.44	1.38	1.43

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	F	2	NAG	O5-C1	-3.02	1.38	1.43

The worst 5 of 102 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	K	1	NAG	C1-O5-C5	6.54	121.05	112.19
3	J	5	MAN	C1-O5-C5	6.36	120.81	112.19
2	K	1	NAG	O5-C1-C2	6.32	121.27	111.29
3	J	6	MAN	O5-C5-C6	5.58	115.94	107.20
2	E	1	NAG	C1-O5-C5	5.16	119.19	112.19

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
2	K	1	NAG	C1

5 of 12 torsion outliers are listed below:

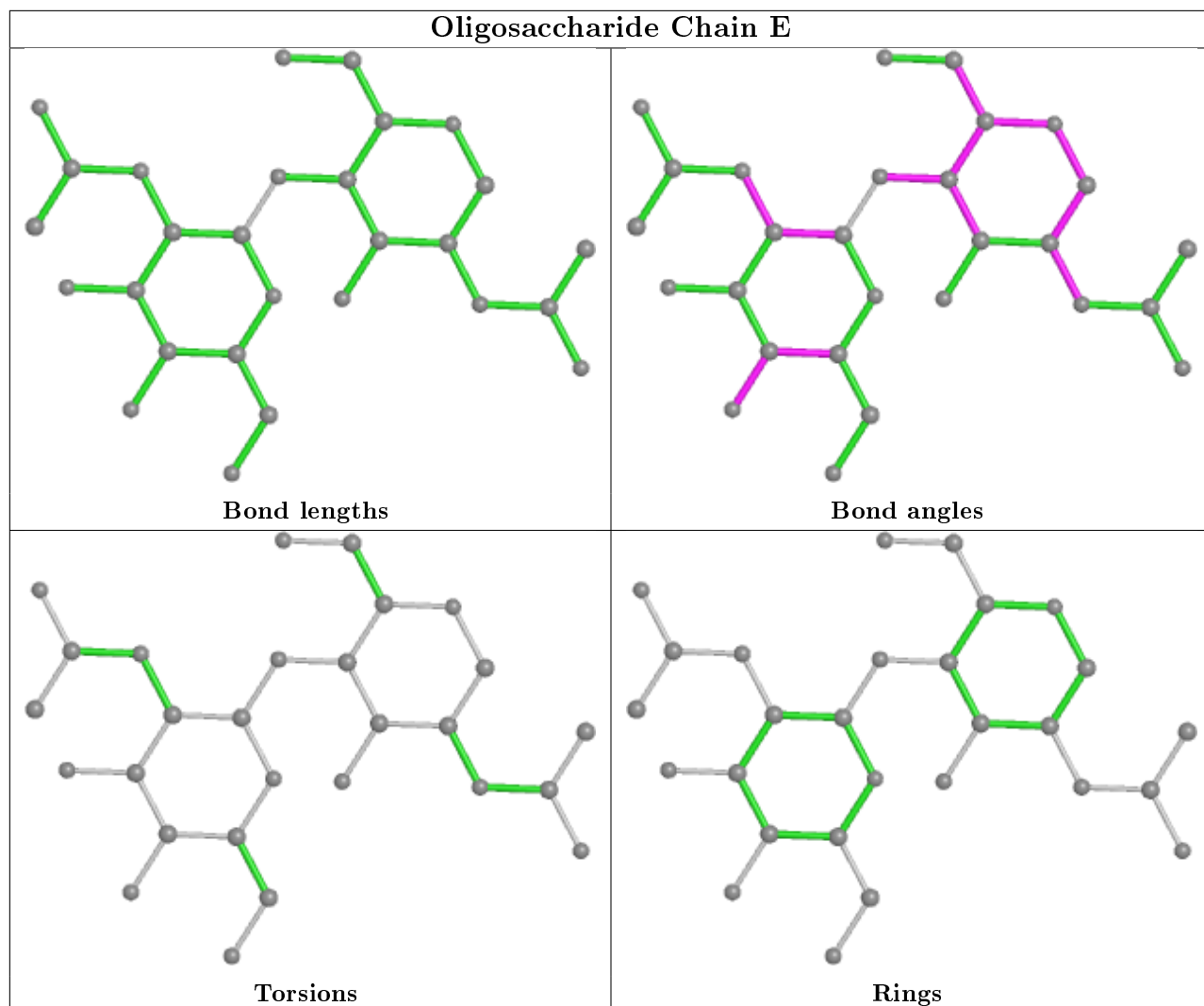
Mol	Chain	Res	Type	Atoms
3	H	6	MAN	O5-C5-C6-O6
3	J	6	MAN	O5-C5-C6-O6
3	L	6	MAN	C4-C5-C6-O6
3	F	6	MAN	C4-C5-C6-O6
3	H	5	MAN	C4-C5-C6-O6

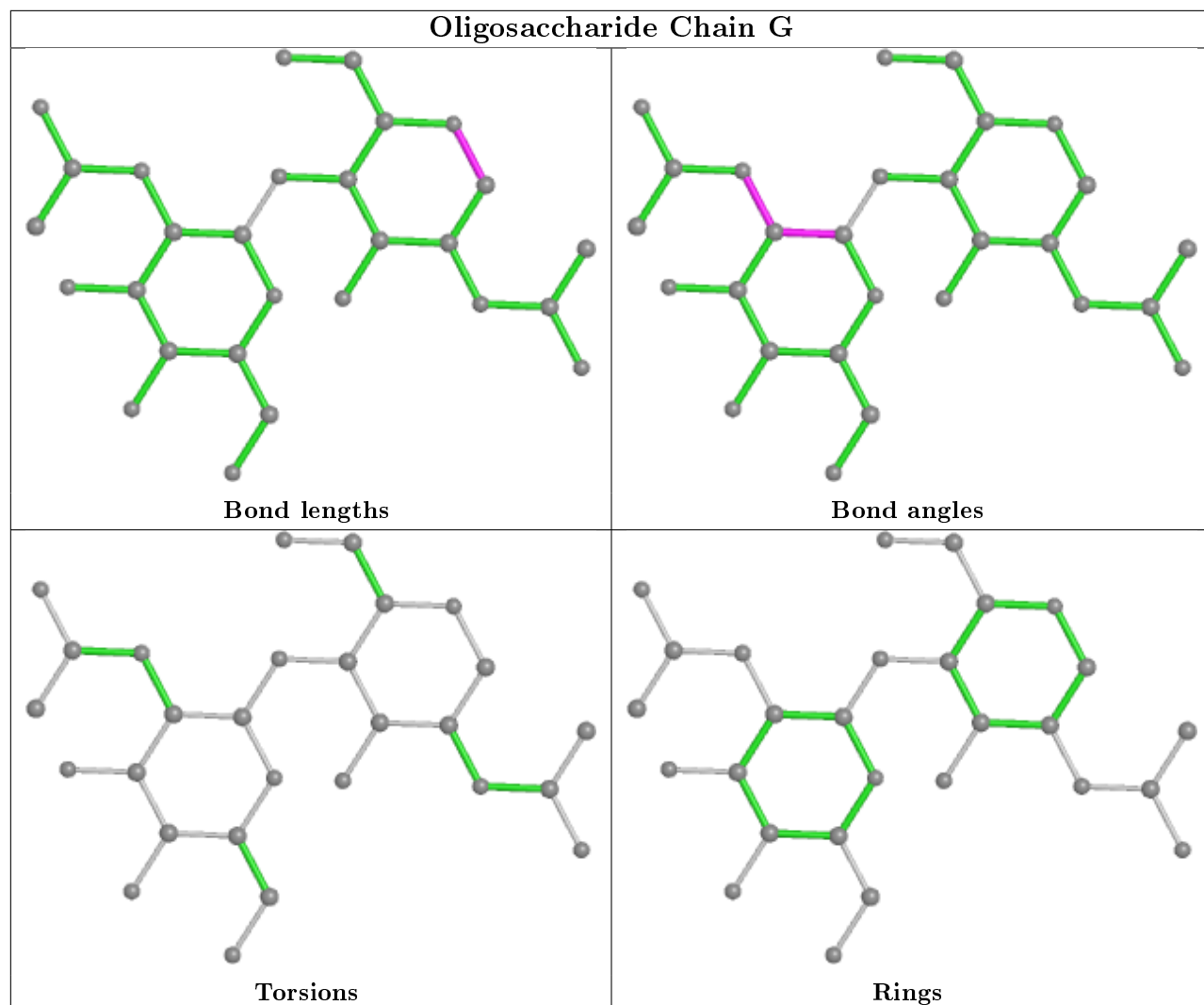
There are no ring outliers.

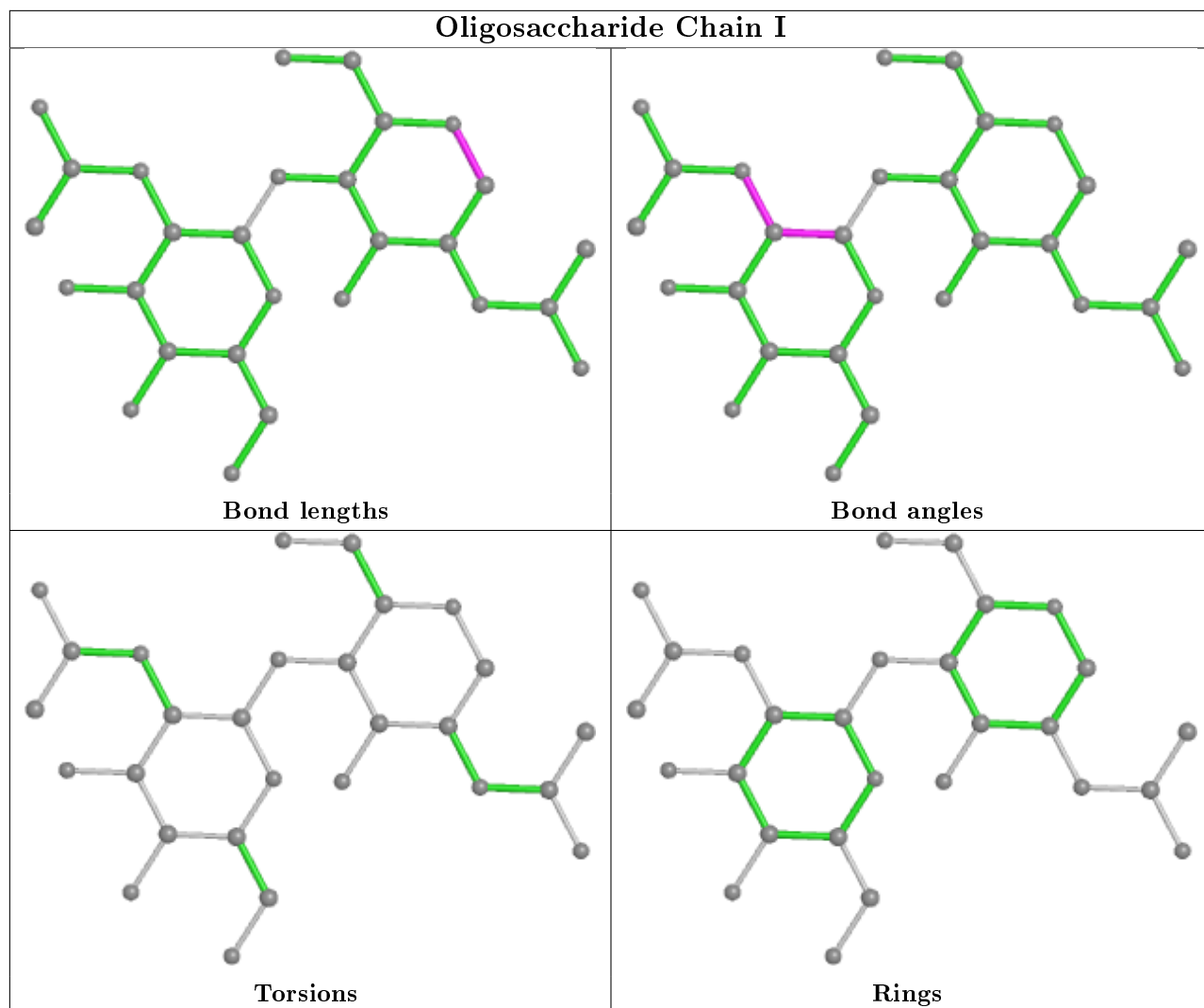
6 monomers are involved in 3 short contacts:

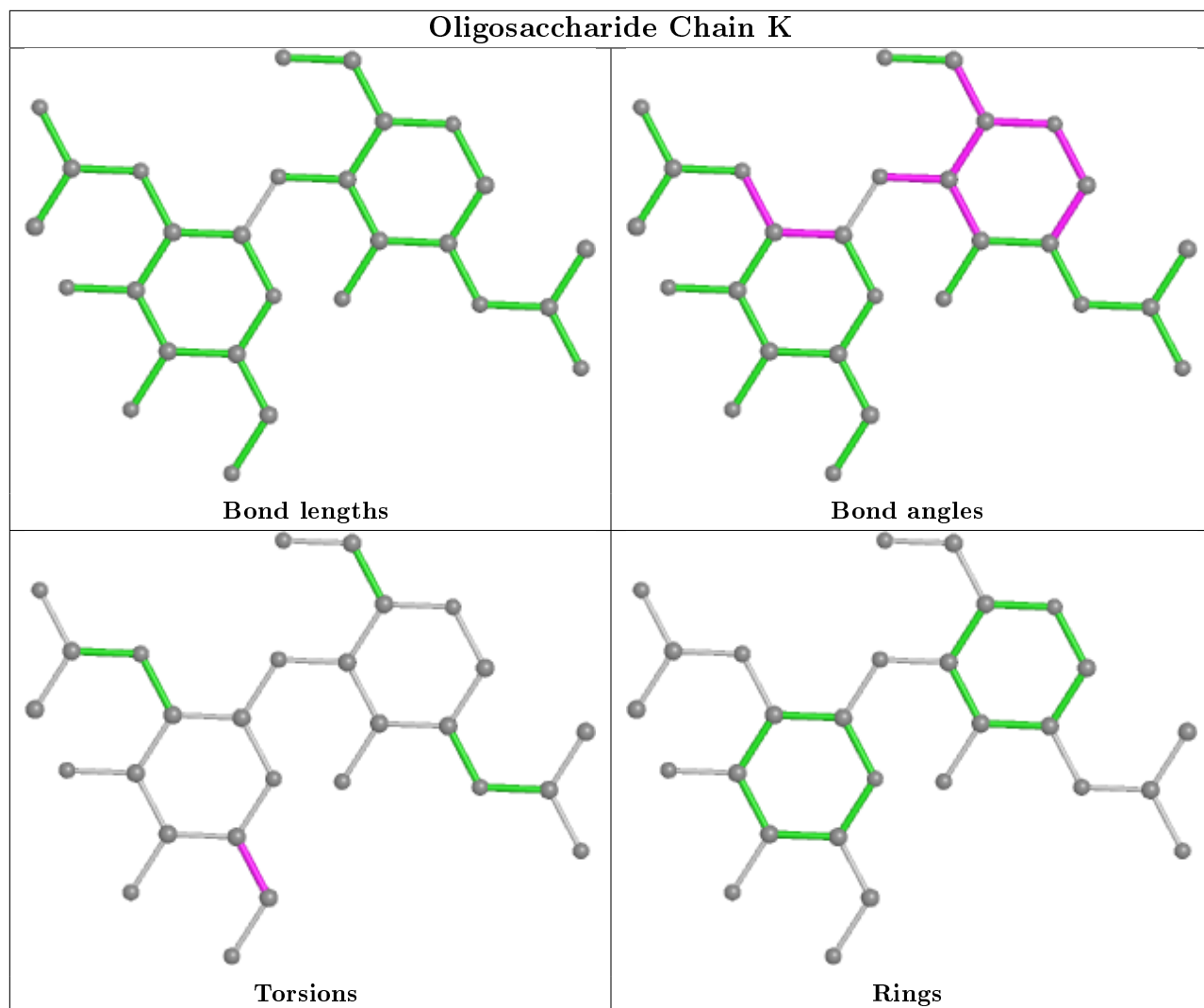
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	J	6	MAN	1	0
2	E	1	NAG	1	0
3	H	3	BMA	1	0
3	H	6	MAN	1	0
2	E	2	NAG	1	0
3	J	3	BMA	1	0

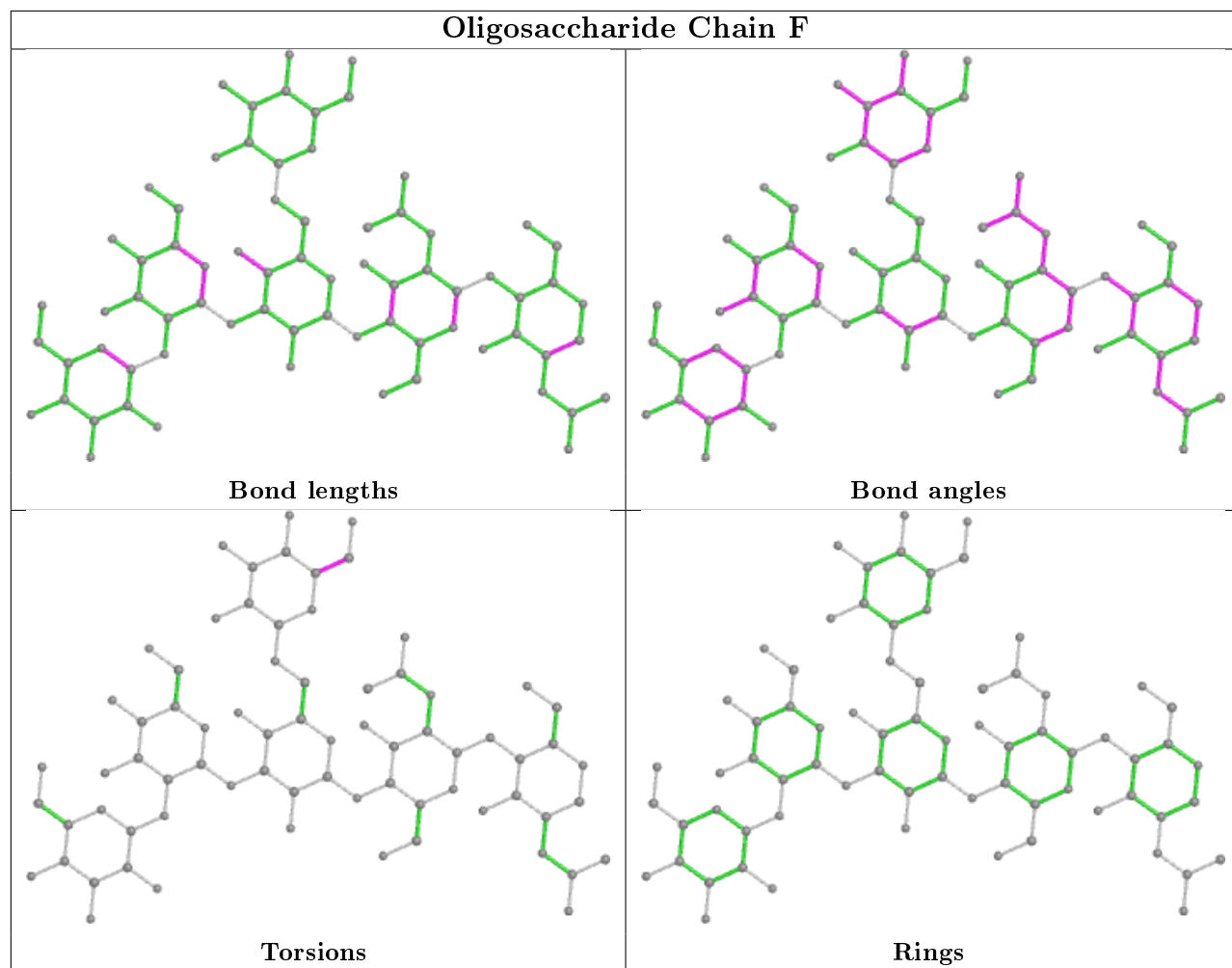
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for oligosaccharide.

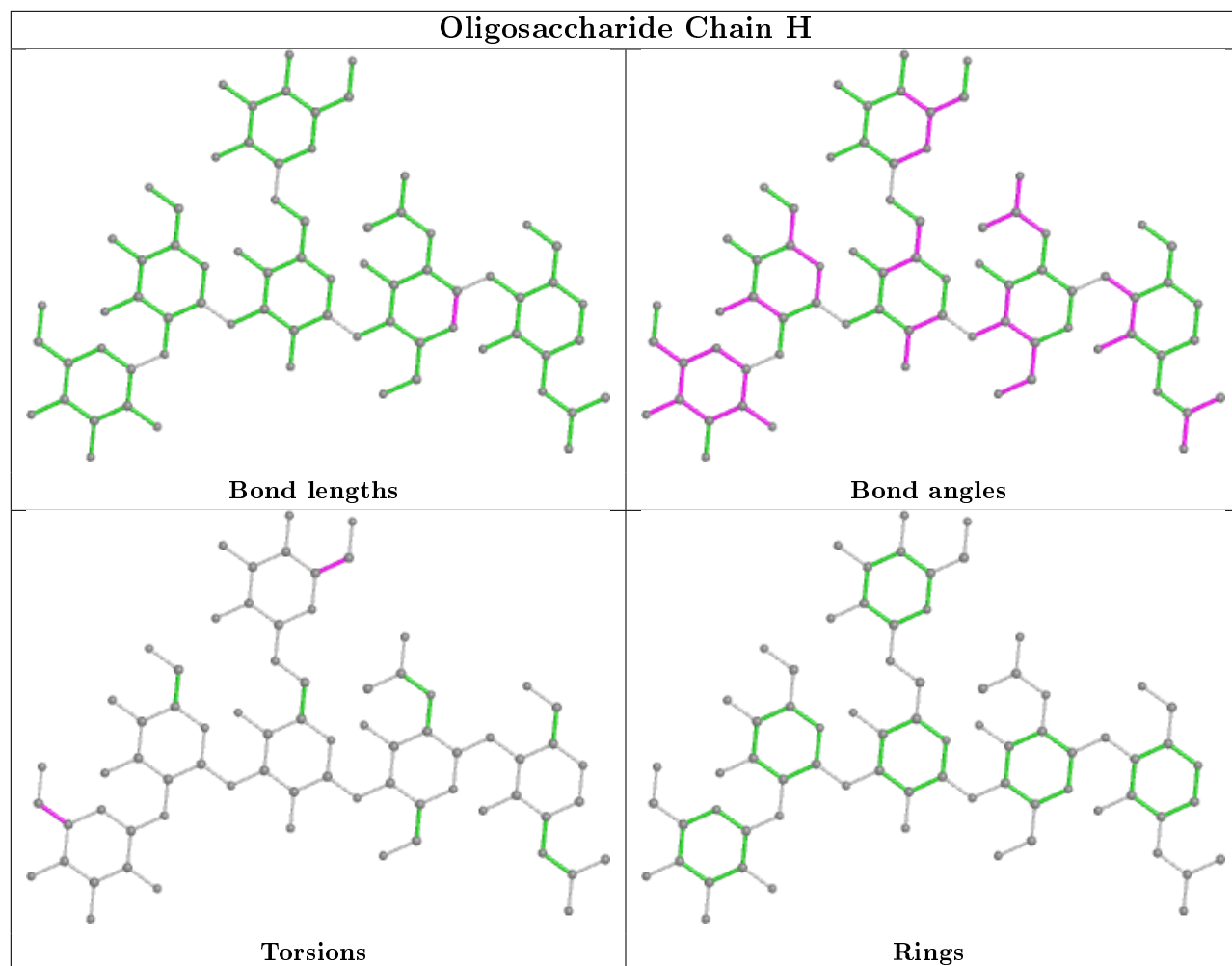


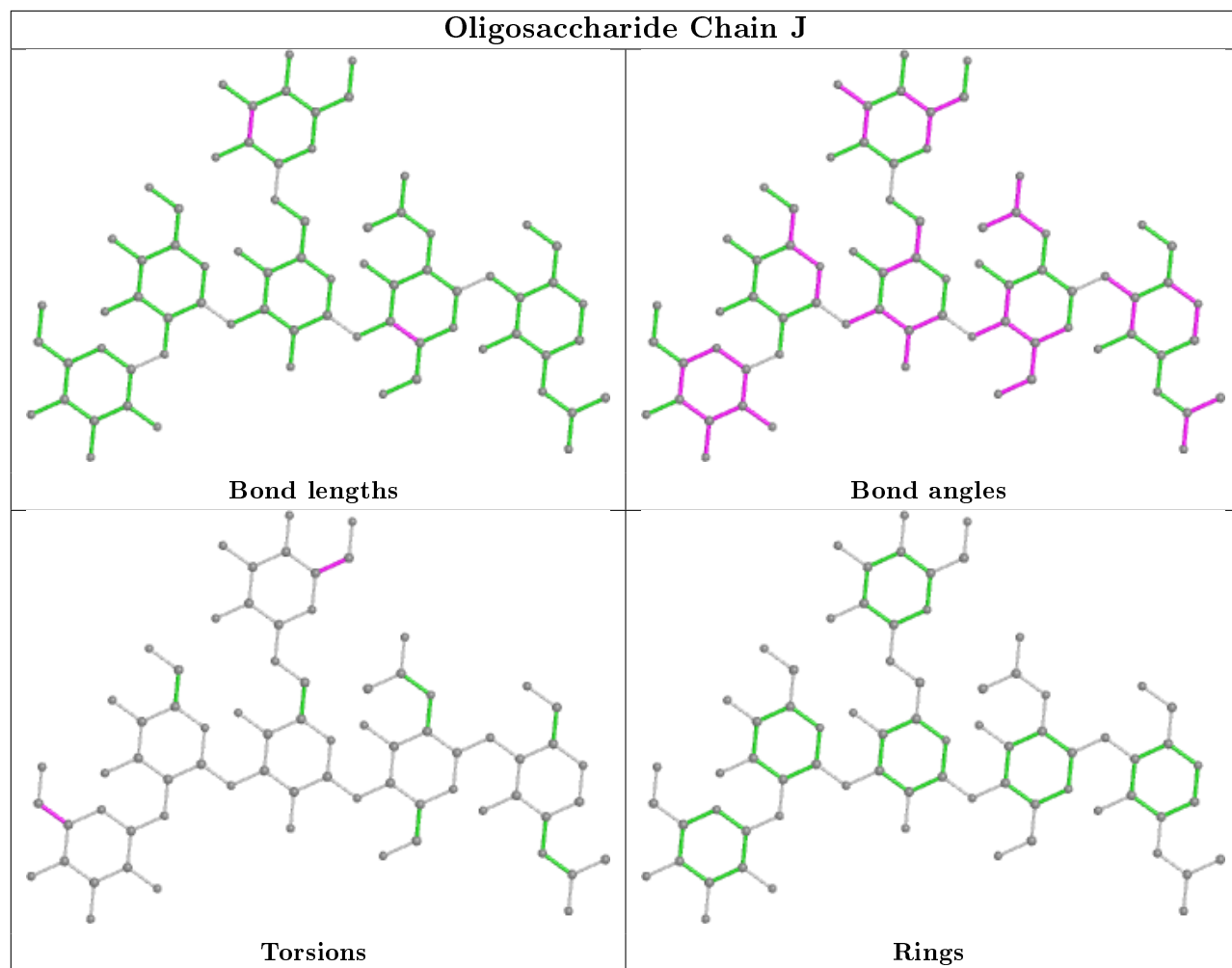


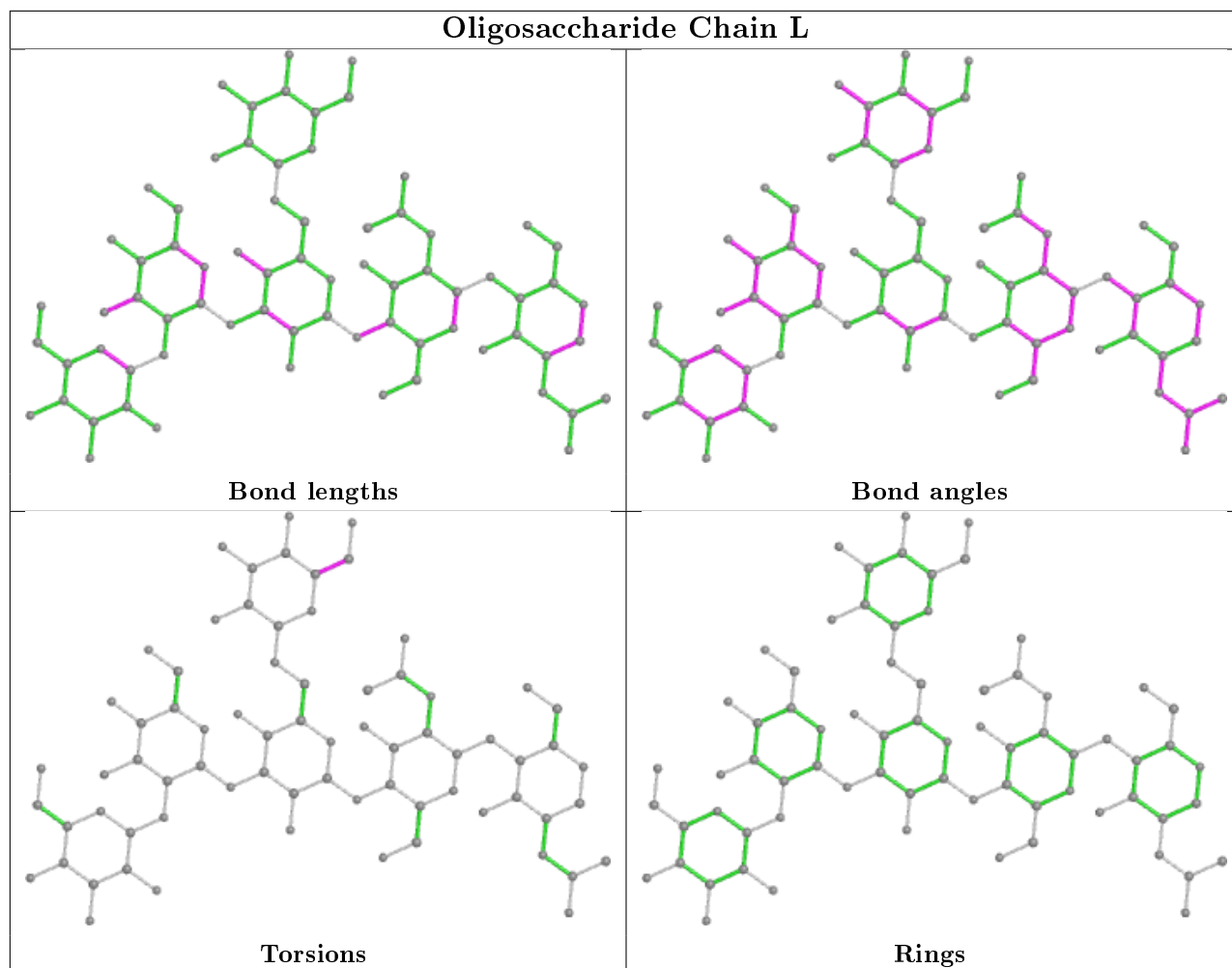












5.6 Ligand geometry [i](#)

Of 8 ligands modelled in this entry, 4 are monoatomic - leaving 4 for Mogul analysis.

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	D	504	1	14,14,15	0.66	0	17,19,21	1.44	3 (17%)
5	NAG	B	504	1	14,14,15	0.51	0	17,19,21	1.25	2 (11%)
5	NAG	C	504	1	14,14,15	0.56	0	17,19,21	1.33	2 (11%)
5	NAG	A	504	1	14,14,15	0.69	0	17,19,21	1.26	1 (5%)

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	D	504	1	-	0/6/23/26	0/1/1/1
5	NAG	B	504	1	-	2/6/23/26	0/1/1/1
5	NAG	C	504	1	-	2/6/23/26	0/1/1/1
5	NAG	A	504	1	-	0/6/23/26	0/1/1/1

There are no bond length outliers.

The worst 5 of 8 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	D	504	NAG	O5-C5-C6	3.45	112.62	107.20
5	A	504	NAG	O5-C5-C6	2.96	111.85	107.20
5	B	504	NAG	O5-C1-C2	2.79	115.69	111.29
5	B	504	NAG	C1-O5-C5	2.73	115.89	112.19
5	D	504	NAG	C3-C4-C5	-2.64	105.52	110.24

There are no chirality outliers.

All (4) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	C	504	NAG	C4-C5-C6-O6
5	B	504	NAG	C4-C5-C6-O6
5	B	504	NAG	O5-C5-C6-O6
5	C	504	NAG	O5-C5-C6-O6

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data [i](#)

6.1 Protein, DNA and RNA chains [i](#)

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	388/397 (97%)	0.01	8 (2%) 63 59	15, 23, 37, 91	0
1	B	388/397 (97%)	0.33	15 (3%) 39 33	21, 29, 43, 124	0
1	C	388/397 (97%)	0.15	8 (2%) 63 59	21, 29, 43, 118	0
1	D	388/397 (97%)	-0.01	5 (1%) 77 74	15, 23, 36, 88	0
All	All	1552/1588 (97%)	0.12	36 (2%) 60 56	15, 27, 40, 124	0

The worst 5 of 36 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	149	ILE	22.4
1	C	149	ILE	14.5
1	A	149	ILE	11.6
1	D	149	ILE	8.9
1	B	150	HIS	6.0

6.2 Non-standard residues in protein, DNA, RNA chains [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
3	MAN	H	6	11/12	0.53	0.34	50,57,62,64	0
5	NAG	B	504	14/15	0.54	0.29	63,70,76,83	0
5	NAG	C	504	14/15	0.56	0.32	59,68,74,83	0
2	NAG	E	2	14/15	0.64	0.31	57,64,69,73	0
3	MAN	J	6	11/12	0.70	0.25	47,52,58,59	0
5	NAG	D	504	14/15	0.72	0.20	57,62,65,67	0
3	MAN	H	5	11/12	0.73	0.25	49,51,55,58	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
5	NAG	A	504	14/15	0.74	0.15	56,61,62,63	0
2	NAG	G	2	14/15	0.74	0.24	43,67,75,80	0
2	NAG	K	2	14/15	0.74	0.28	53,65,70,75	0
3	MAN	J	5	11/12	0.75	0.22	49,53,56,61	0
2	NAG	I	1	14/15	0.76	0.19	43,48,55,58	0
2	NAG	K	1	14/15	0.79	0.19	43,48,56,59	0
2	NAG	E	1	14/15	0.80	0.16	44,51,59,62	0
3	MAN	L	5	11/12	0.81	0.15	40,49,54,55	0
2	NAG	I	2	14/15	0.81	0.21	46,63,67,71	0
3	MAN	L	6	11/12	0.82	0.19	41,45,50,52	0
3	MAN	F	6	11/12	0.84	0.17	40,43,49,50	0
2	NAG	G	1	14/15	0.85	0.15	43,48,55,60	0
3	MAN	F	5	11/12	0.88	0.13	40,46,52,56	0
3	MAN	H	4	11/12	0.89	0.12	32,34,38,42	0
3	NAG	J	1	14/15	0.89	0.12	26,33,41,42	0
3	NAG	J	2	14/15	0.90	0.12	26,30,34,35	0
3	NAG	H	1	14/15	0.91	0.11	26,33,40,41	0
3	MAN	J	4	11/12	0.91	0.12	31,32,35,40	0
3	NAG	L	1	14/15	0.91	0.14	21,24,38,40	0
3	NAG	H	2	14/15	0.93	0.10	27,30,32,34	0
3	MAN	L	4	11/12	0.94	0.10	27,29,33,36	0
3	NAG	F	2	14/15	0.95	0.10	21,22,23,24	0
3	MAN	F	4	11/12	0.95	0.10	28,29,32,36	0
3	NAG	L	2	14/15	0.95	0.08	22,23,24,24	0
3	NAG	F	1	14/15	0.96	0.10	22,24,37,42	0

6.3 Carbohydrates [\(i\)](#)

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
3	MAN	H	6	11/12	0.53	0.34	50,57,62,64	0
2	NAG	E	2	14/15	0.64	0.31	57,64,69,73	0
3	MAN	J	6	11/12	0.70	0.25	47,52,58,59	0
3	MAN	H	5	11/12	0.73	0.25	49,51,55,58	0
2	NAG	G	2	14/15	0.74	0.24	43,67,75,80	0
2	NAG	K	2	14/15	0.74	0.28	53,65,70,75	0
3	MAN	J	5	11/12	0.75	0.22	49,53,56,61	0
2	NAG	I	1	14/15	0.76	0.19	43,48,55,58	0

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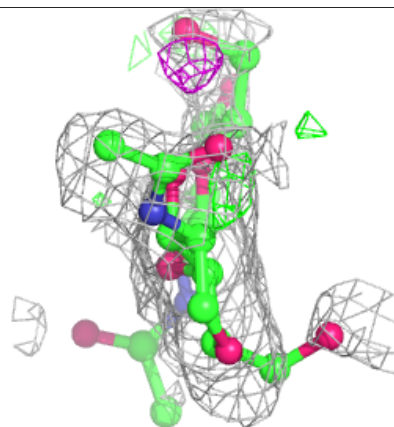
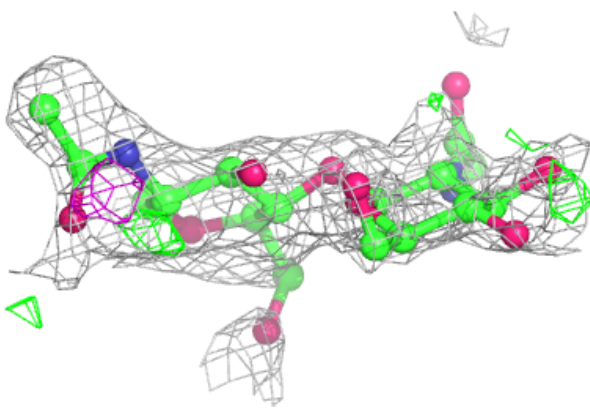
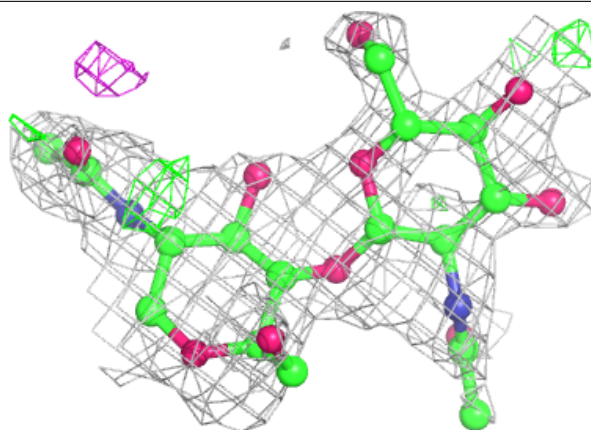
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
2	NAG	K	1	14/15	0.79	0.19	43,48,56,59	0
2	NAG	E	1	14/15	0.80	0.16	44,51,59,62	0
2	NAG	I	2	14/15	0.81	0.21	46,63,67,71	0
3	MAN	L	5	11/12	0.81	0.15	40,49,54,55	0
3	MAN	L	6	11/12	0.82	0.19	41,45,50,52	0
3	MAN	F	6	11/12	0.84	0.17	40,43,49,50	0
2	NAG	G	1	14/15	0.85	0.15	43,48,55,60	0
3	MAN	F	5	11/12	0.88	0.13	40,46,52,56	0
3	NAG	J	1	14/15	0.89	0.12	26,33,41,42	0
3	MAN	H	4	11/12	0.89	0.12	32,34,38,42	0
3	NAG	J	2	14/15	0.90	0.12	26,30,34,35	0
3	NAG	L	1	14/15	0.91	0.14	21,24,38,40	0
3	NAG	H	1	14/15	0.91	0.11	26,33,40,41	0
3	MAN	J	4	11/12	0.91	0.12	31,32,35,40	0
3	BMA	H	3	11/12	0.91	0.09	27,30,37,40	0
3	BMA	J	3	11/12	0.92	0.13	26,31,37,37	0
3	BMA	L	3	11/12	0.92	0.12	22,25,30,33	0
3	NAG	H	2	14/15	0.93	0.10	27,30,32,34	0
3	MAN	L	4	11/12	0.94	0.10	27,29,33,36	0
3	NAG	L	2	14/15	0.95	0.08	22,23,24,24	0
3	MAN	F	4	11/12	0.95	0.10	28,29,32,36	0
3	NAG	F	2	14/15	0.95	0.10	21,22,23,24	0
3	BMA	F	3	11/12	0.96	0.08	21,25,28,31	0
3	NAG	F	1	14/15	0.96	0.10	22,24,37,42	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.

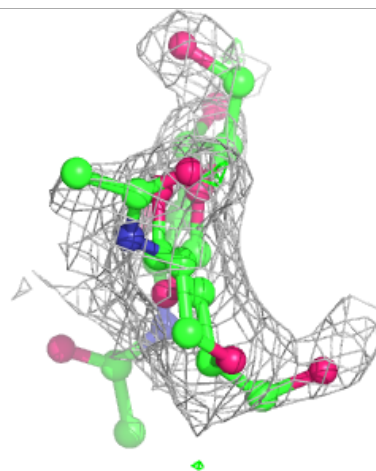
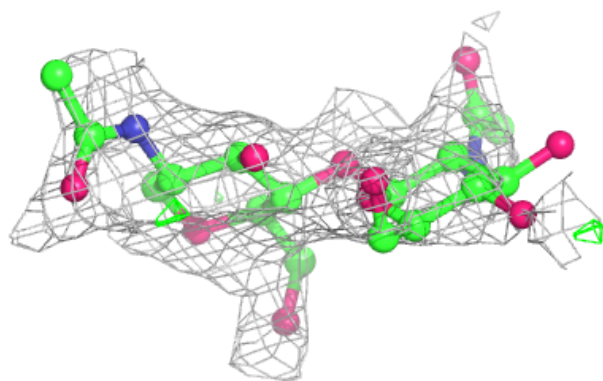
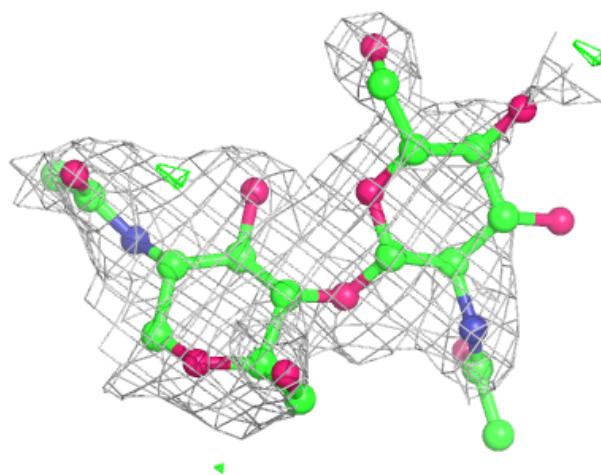
Electron density around Chain E:

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



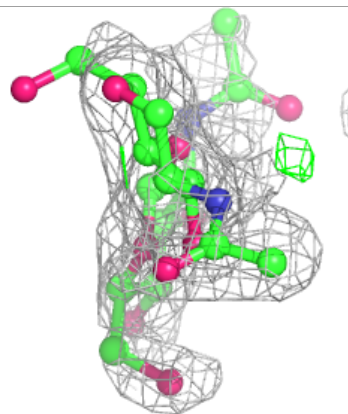
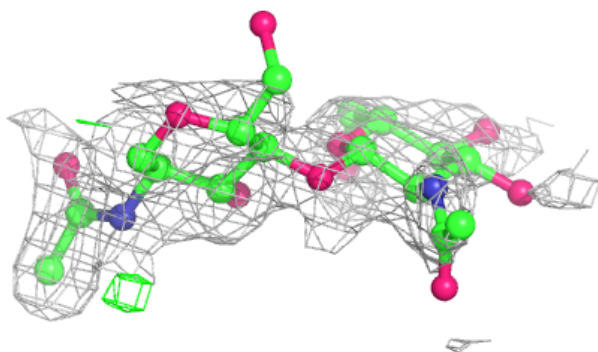
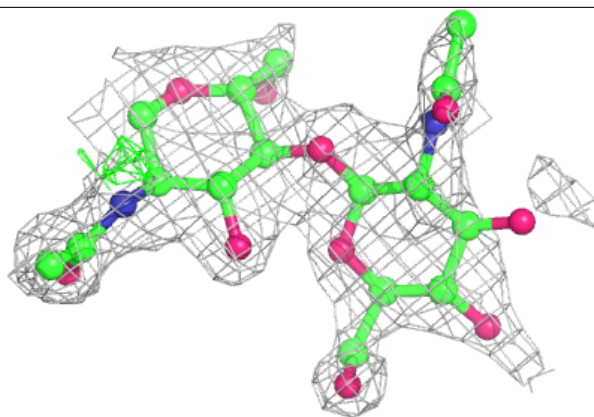
Electron density around Chain G:

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

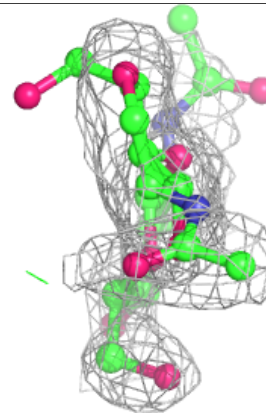
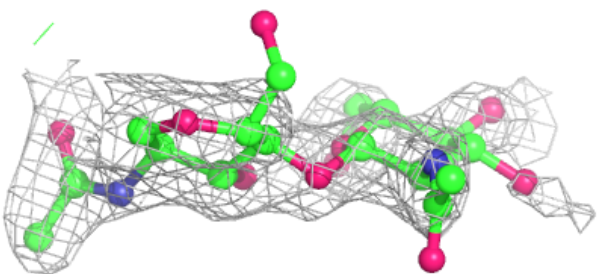
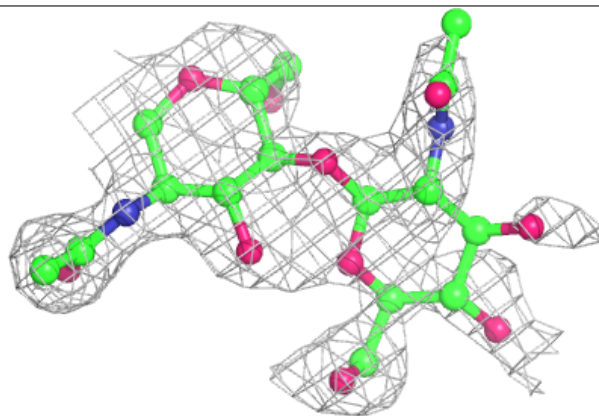


Electron density around Chain I:

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

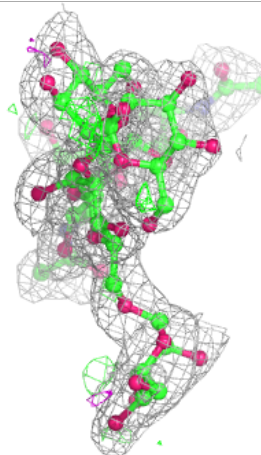
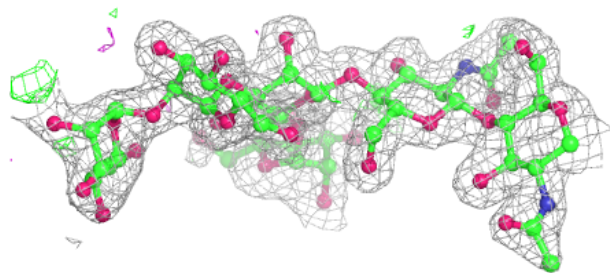
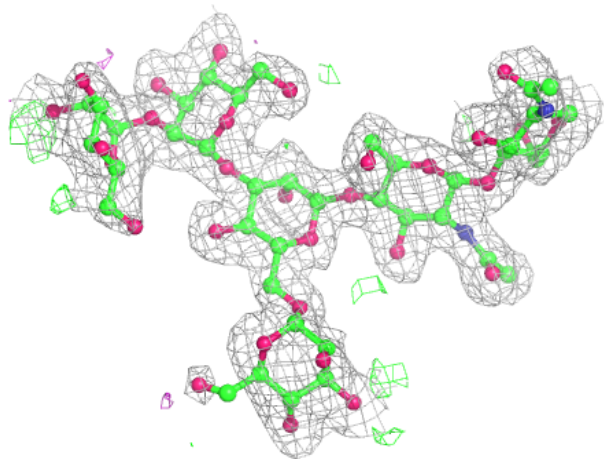
**Electron density around Chain K:**

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



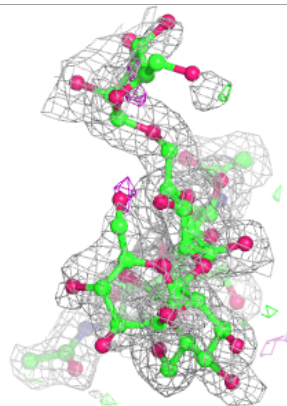
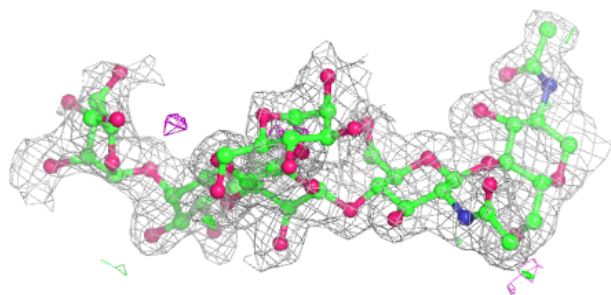
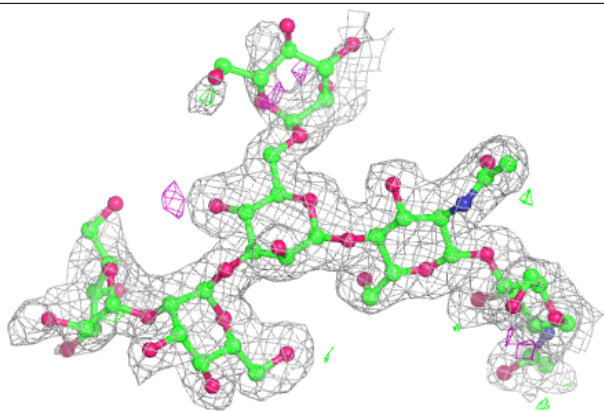
Electron density around Chain F:

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

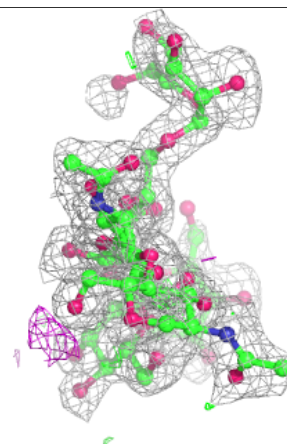
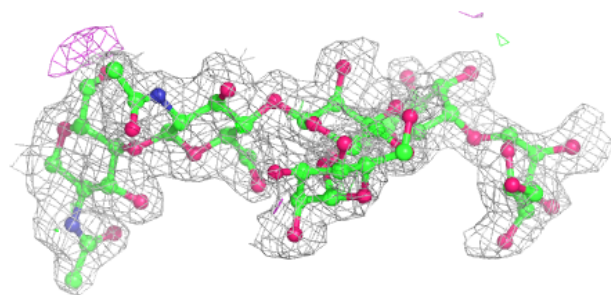
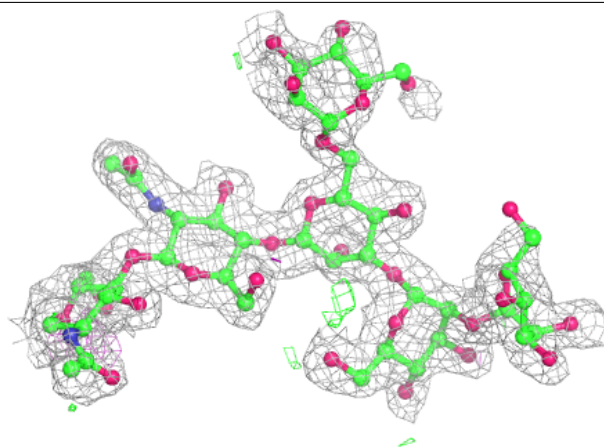


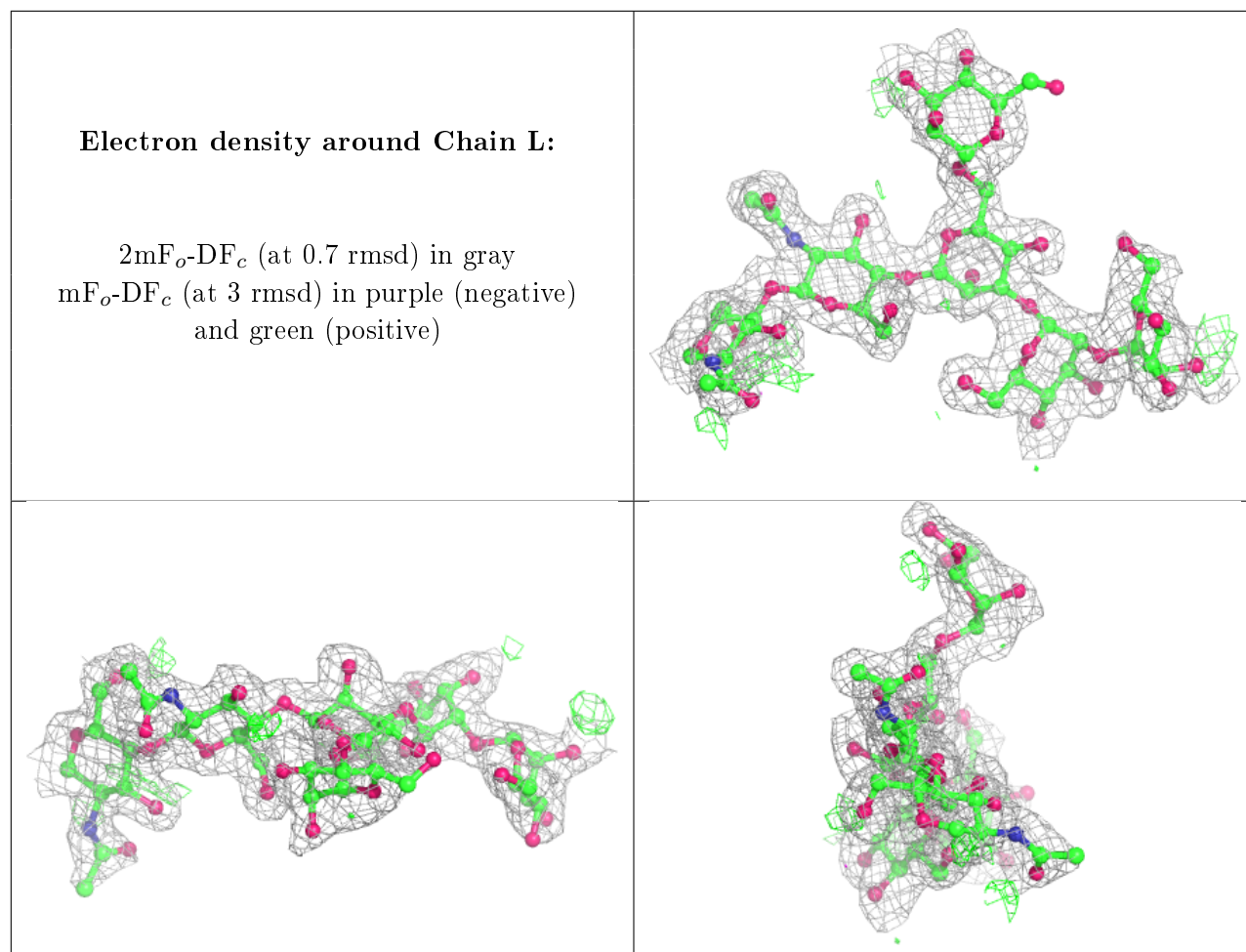
Electron density around Chain H:

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

**Electron density around Chain J:**

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





6.4 Ligands [i](#)

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
5	NAG	B	504	14/15	0.54	0.29	63,70,76,83	0
5	NAG	C	504	14/15	0.56	0.32	59,68,74,83	0
5	NAG	D	504	14/15	0.72	0.20	57,62,65,67	0
5	NAG	A	504	14/15	0.74	0.15	56,61,62,63	0
4	CA	A	501	1/1	0.99	0.04	24,24,24,24	0
4	CA	D	501	1/1	0.99	0.05	24,24,24,24	0
4	CA	B	501	1/1	0.99	0.07	26,26,26,26	0
4	CA	C	501	1/1	1.00	0.06	26,26,26,26	0

6.5 Other polymers

There are no such residues in this entry.