



wwPDB X-ray Structure Validation Summary Report ⓘ

Aug 7, 2020 – 10:48 PM BST

PDB ID : 4QZV
Title : Bat-derived coronavirus HKU4 uses MERS-CoV receptor human CD26 for cell entry
Authors : Gao, F.G.; Wang, Q.H.; Qi, J.X.; Lu, G.W.
Deposited on : 2014-07-29
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 ⓘ 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

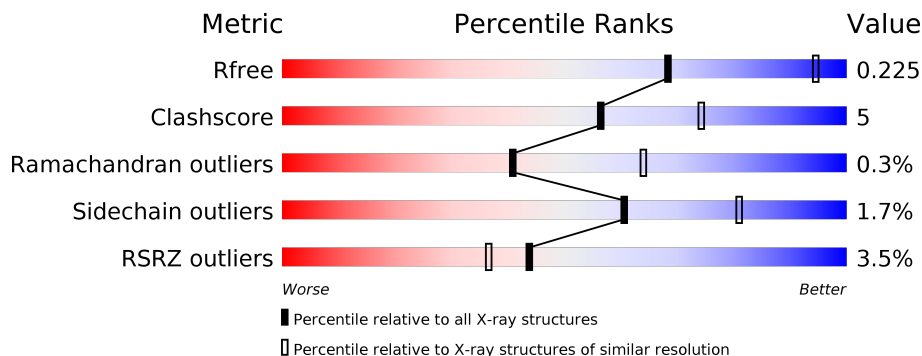
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 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 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	734	
1	C	734	
2	B	246	
2	D	246	
3	E	2	
3	G	2	

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Mol	Chain	Length	Quality of chain
3	H	2	 50% 50%
3	I	2	 50% 50%
3	K	2	 100%
3	L	2	 50% 50%
4	F	3	 33% 67%
4	J	3	 33% 67%

2 Entry composition [i](#)

There are 6 unique types of molecules in this entry. The entry contains 15900 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 Dipeptidyl peptidase 4.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	727	Total	C	N	O	S	0	0	0
			5957	3824	981	1126	26			
1	C	727	Total	C	N	O	S	0	0	0
			5957	3824	981	1126	26			

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	767	HIS	-	expression tag	UNP P27487
A	768	HIS	-	expression tag	UNP P27487
A	769	HIS	-	expression tag	UNP P27487
A	770	HIS	-	expression tag	UNP P27487
A	771	HIS	-	expression tag	UNP P27487
A	772	HIS	-	expression tag	UNP P27487
C	767	HIS	-	expression tag	UNP P27487
C	768	HIS	-	expression tag	UNP P27487
C	769	HIS	-	expression tag	UNP P27487
C	770	HIS	-	expression tag	UNP P27487
C	771	HIS	-	expression tag	UNP P27487
C	772	HIS	-	expression tag	UNP P27487

- Molecule 2 is a protein called Spike protein S1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	B	208	Total	C	N	O	S	0	0	0
			1597	1017	260	306	14			
2	D	208	Total	C	N	O	S	0	0	0
			1597	1017	260	306	14			

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	241	HIS	-	expression tag	UNP A3EX94
B	242	HIS	-	expression tag	UNP A3EX94
B	243	HIS	-	expression tag	UNP A3EX94
B	244	HIS	-	expression tag	UNP A3EX94
B	245	HIS	-	expression tag	UNP A3EX94
B	246	HIS	-	expression tag	UNP A3EX94
D	241	HIS	-	expression tag	UNP A3EX94
D	242	HIS	-	expression tag	UNP A3EX94
D	243	HIS	-	expression tag	UNP A3EX94
D	244	HIS	-	expression tag	UNP A3EX94
D	245	HIS	-	expression tag	UNP A3EX94
D	246	HIS	-	expression tag	UNP A3EX94

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



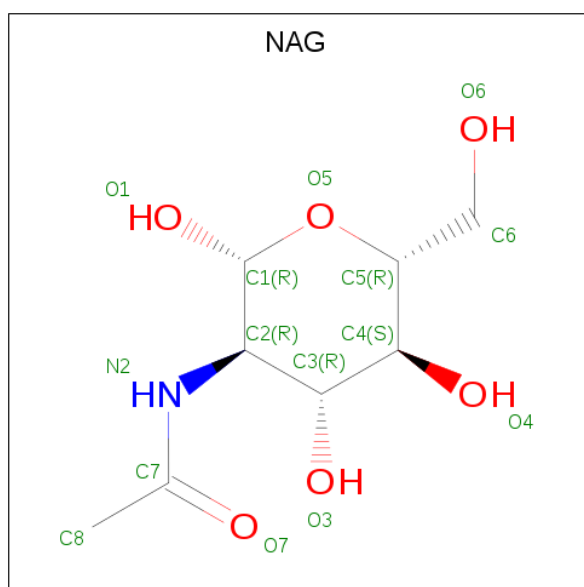
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
			Total	C	N	O			
3	E	2	28	16	2	10	0	0	0
3	G	2	28	16	2	10	0	0	0
3	H	2	28	16	2	10	0	0	0
3	I	2	28	16	2	10	0	0	0
3	K	2	28	16	2	10	0	0	0
3	L	2	28	16	2	10	0	0	0

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



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
4	F	3	Total	C	N	O	0	0	0
			39	22	2	15			
4	J	3	Total	C	N	O	0	0	0
			39	22	2	15			

- 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	0	0
			14	8	1	5		
5	A	1	Total	C	N	O	0	0
			14	8	1	5		
5	A	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	C	1	Total	C	N	O	0	0
			14	8	1	5		

- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	195	Total	O	0	0
			195	195		
6	B	26	Total	O	0	0
			26	26		

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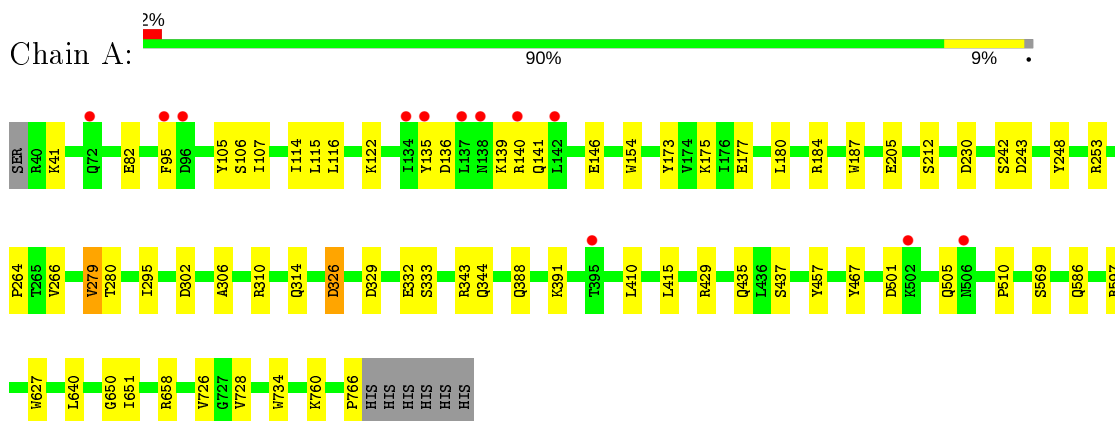
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	C	239	Total 239	O 239	0	0
6	D	16	Total 16	O 16	0	0

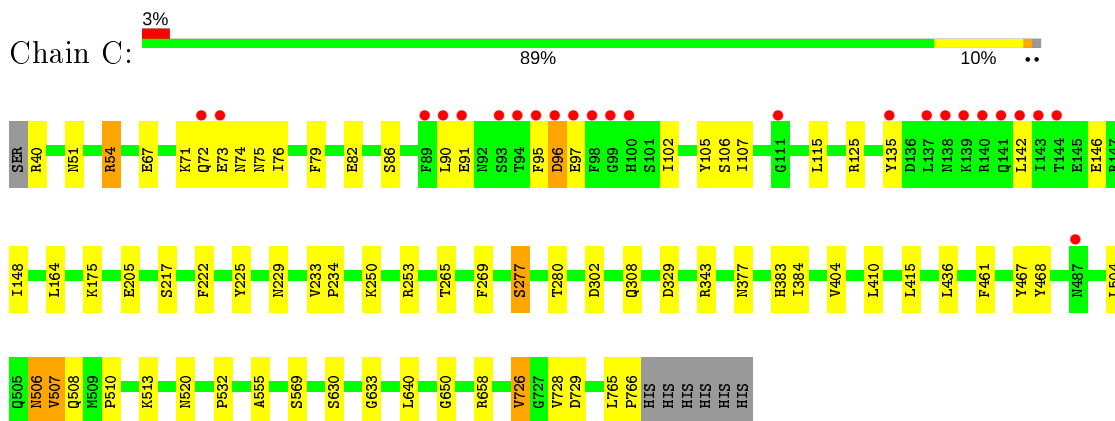
3 Residue-property plots [i](#)

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.

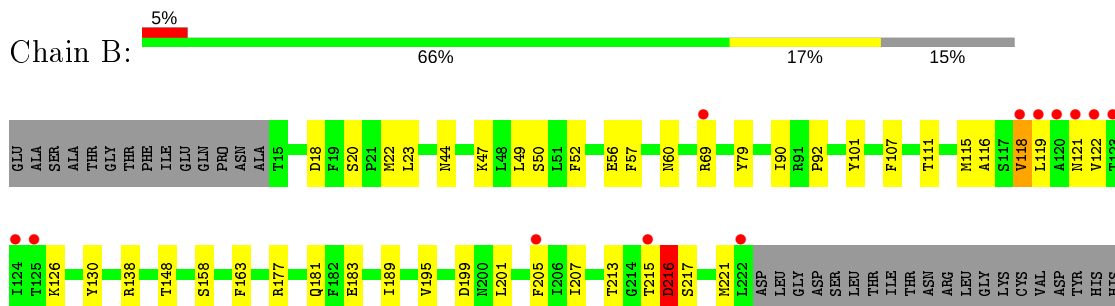
- Molecule 1: Dipeptidyl peptidase 4



- Molecule 1: Dipeptidyl peptidase 4



- Molecule 2: Spike protein S1



HIS
HIS
HIS
HIS

- Molecule 2: Spike protein S1

Chain D:  7% 65% 16% 15%

GLU ALA SER ALA THR GLY THR PHE ILE GLU GLN PRO ASN ALA T15 M22 M44 L49 F52 F57 S58 C59 S63 P64 D65 S66 I67 A68 R69 Y72 S73 Y79 F80 P83 M86 K87 S88 S94 A95 G96 N97 I98 P99 L100 Y101 M102 A116 S117

V118 L119 A120 M121 V122 T123 I124 T125 K126 R138 I150 Y151 Y157 R161 D162 E170 R177 T180 Q181 F182 E183 G184 T185 T193 D199 S204 Y211 G212 G214 T215 D216 S217 V218 M221 L222 ASP LEU LEU LEU SER LEU THR ILE THR ASN ARG LEU

GLY LYS CYS VAL ASP TYR HIS HIS HIS HIS HIS HIS

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

Chain E:  100%

MAG1
MAG2

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

Chain G:  50% 50%

MAG1
MAG2

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

Chain H:  50% 50%

MAG1
MAG2

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

Chain I:  50% 50%

MAG1
MAG2

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

Chain K:  100%


MAC1
MAC2

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

Chain L:  50% 50%

MAC1
MAC2

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

Chain F:  33% 67%

MAC1
MAC2
BMA3

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

Chain J:  33% 67%

MAC1
MAC2
BMA3

4 Data and refinement statistics

Property	Value	Source
Space group	P 2 21 21	Depositor
Cell constants a, b, c, α , β , γ	78.80Å 203.17Å 207.75Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	46.30 – 2.59 46.30 – 2.59	Depositor EDS
% Data completeness (in resolution range)	98.0 (46.30-2.59) 93.2 (46.30-2.59)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.49 (at 2.58Å)	Xtrriage
Refinement program	PHENIX (phenix.refine: 1.8.4_1496)	Depositor
R, R_{free}	0.190 , 0.224 0.193 , 0.225	Depositor DCC
R_{free} test set	5113 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	36.7	Xtrriage
Anisotropy	0.732	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 37.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtrriage
Estimated twinning fraction	0.007 for -h,l,k	Xtrriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	15900	wwPDB-VP
Average B, all atoms (Å ²)	47.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.64% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: BMA, 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	0.29	0/6129	0.46	0/8336
1	C	0.32	0/6129	0.49	1/8336 (0.0%)
2	B	0.39	0/1635	0.49	0/2222
2	D	0.40	0/1635	0.51	0/2222
All	All	0.32	0/15528	0.48	1/21116 (0.0%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	C	0	1
2	B	0	1
2	D	0	2
All	All	0	4

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	96	ASP	CB-CG-OD2	5.28	123.05	118.30

There are no chirality outliers.

All (4) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	B	121	ASN	Peptide
1	C	96	ASP	Peptide

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Mol	Chain	Res	Type	Group
2	D	121	ASN	Peptide
2	D	215	THR	Peptide

5.2 Too-close contacts [i](#)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	5957	0	5673	45	0
1	C	5957	0	5674	54	0
2	B	1597	0	1553	26	0
2	D	1597	0	1553	28	0
3	E	28	0	25	1	0
3	G	28	0	25	1	0
3	H	28	0	25	0	0
3	I	28	0	25	0	0
3	K	28	0	25	0	0
3	L	28	0	25	0	0
4	F	39	0	34	0	0
4	J	39	0	34	2	0
5	A	42	0	39	0	0
5	C	28	0	26	0	0
6	A	195	0	0	12	0
6	B	26	0	0	1	0
6	C	239	0	0	12	0
6	D	16	0	0	2	0
All	All	15900	0	14736	152	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 5.

The worst 5 of 152 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:A:658:ARG:NH1	6:A:1029:HOH:O	1.88	1.06
1:A:253:ARG:HH21	1:C:253:ARG:HH21	1.16	0.91
2:D:180:THR:HG22	2:D:183:GLU:HG2	1.59	0.84

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:63:SER:OG	2:D:66:SER:OG	1.96	0.84
2:D:138:ARG:NH2	2:D:183:GLU:O	2.11	0.83

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	725/734 (99%)	694 (96%)	31 (4%)	0	100	100
1	C	725/734 (99%)	688 (95%)	34 (5%)	3 (0%)	34	57
2	B	206/246 (84%)	189 (92%)	16 (8%)	1 (0%)	29	52
2	D	206/246 (84%)	195 (95%)	10 (5%)	1 (0%)	29	52
All	All	1862/1960 (95%)	1766 (95%)	91 (5%)	5 (0%)	41	64

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	216	ASP
1	C	97	GLU
1	C	520	ASN
1	C	73	GLU
2	D	216	ASP

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	652/659 (99%)	648 (99%)	4 (1%)	86	95
1	C	652/659 (99%)	644 (99%)	8 (1%)	71	87
2	B	180/212 (85%)	174 (97%)	6 (3%)	38	64
2	D	180/212 (85%)	169 (94%)	11 (6%)	18	38
All	All	1664/1742 (96%)	1635 (98%)	29 (2%)	60	81

5 of 29 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	C	250	LYS
1	C	507	VAL
2	D	161	ARG
1	C	277	SER
1	C	726	VAL

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	731	GLN
2	B	121	ASN
1	C	506	ASN
2	D	153	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](#)

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

5.5 Carbohydrates [i](#)

18 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	E	1	1,3	14,14,15	0.55	0	17,19,21	0.86	1 (5%)
3	NAG	E	2	3	14,14,15	0.51	0	17,19,21	0.49	0
4	NAG	F	1	1,4	14,14,15	0.86	1 (7%)	17,19,21	0.58	0
4	NAG	F	2	4	14,14,15	0.36	0	17,19,21	0.70	0
4	BMA	F	3	4	11,11,12	1.87	3 (27%)	15,15,17	1.72	3 (20%)
3	NAG	G	1	1,3	14,14,15	0.33	0	17,19,21	0.53	0
3	NAG	G	2	3	14,14,15	0.19	0	17,19,21	0.63	0
3	NAG	H	1	1,3	14,14,15	0.63	1 (7%)	17,19,21	0.49	0
3	NAG	H	2	3	14,14,15	0.30	0	17,19,21	0.45	0
3	NAG	I	1	1,3	14,14,15	0.56	0	17,19,21	0.83	1 (5%)
3	NAG	I	2	3	14,14,15	0.38	0	17,19,21	0.35	0
4	NAG	J	1	1,4	14,14,15	0.82	1 (7%)	17,19,21	0.57	0
4	NAG	J	2	4	14,14,15	0.67	1 (7%)	17,19,21	0.42	0
4	BMA	J	3	4	11,11,12	1.86	3 (27%)	15,15,17	1.89	5 (33%)
3	NAG	K	1	1,3	14,14,15	0.27	0	17,19,21	0.52	0
3	NAG	K	2	3	14,14,15	0.25	0	17,19,21	0.62	0
3	NAG	L	1	1,3	14,14,15	0.60	1 (7%)	17,19,21	0.47	0
3	NAG	L	2	3	14,14,15	0.15	0	17,19,21	0.72	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	E	1	1,3	-	0/6/23/26	0/1/1/1
3	NAG	E	2	3	-	2/6/23/26	0/1/1/1
4	NAG	F	1	1,4	-	0/6/23/26	0/1/1/1
4	NAG	F	2	4	-	4/6/23/26	0/1/1/1
4	BMA	F	3	4	-	0/2/19/22	0/1/1/1
3	NAG	G	1	1,3	-	0/6/23/26	0/1/1/1
3	NAG	G	2	3	-	2/6/23/26	0/1/1/1
3	NAG	H	1	1,3	-	0/6/23/26	0/1/1/1
3	NAG	H	2	3	-	4/6/23/26	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	NAG	I	1	1,3	-	0/6/23/26	0/1/1/1
3	NAG	I	2	3	-	2/6/23/26	0/1/1/1
4	NAG	J	1	1,4	-	0/6/23/26	0/1/1/1
4	NAG	J	2	4	-	4/6/23/26	0/1/1/1
4	BMA	J	3	4	-	2/2/19/22	0/1/1/1
3	NAG	K	1	1,3	-	2/6/23/26	0/1/1/1
3	NAG	K	2	3	-	2/6/23/26	0/1/1/1
3	NAG	L	1	1,3	-	0/6/23/26	0/1/1/1
3	NAG	L	2	3	-	4/6/23/26	0/1/1/1

The worst 5 of 11 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	J	3	BMA	O5-C5	3.40	1.50	1.43
4	F	3	BMA	C4-C3	-3.27	1.44	1.52
4	J	3	BMA	C4-C3	-3.17	1.44	1.52
4	F	3	BMA	O5-C5	3.09	1.49	1.43
4	F	3	BMA	C2-C3	-2.99	1.48	1.52

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

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	J	3	BMA	C3-C4-C5	4.68	118.59	110.24
4	F	3	BMA	C3-C4-C5	4.38	118.05	110.24
4	J	3	BMA	C2-C3-C4	2.89	115.90	110.89
3	I	1	NAG	C1-O5-C5	2.78	115.96	112.19
3	E	1	NAG	C1-O5-C5	2.75	115.91	112.19

There are no chirality outliers.

5 of 28 torsion outliers are listed below:

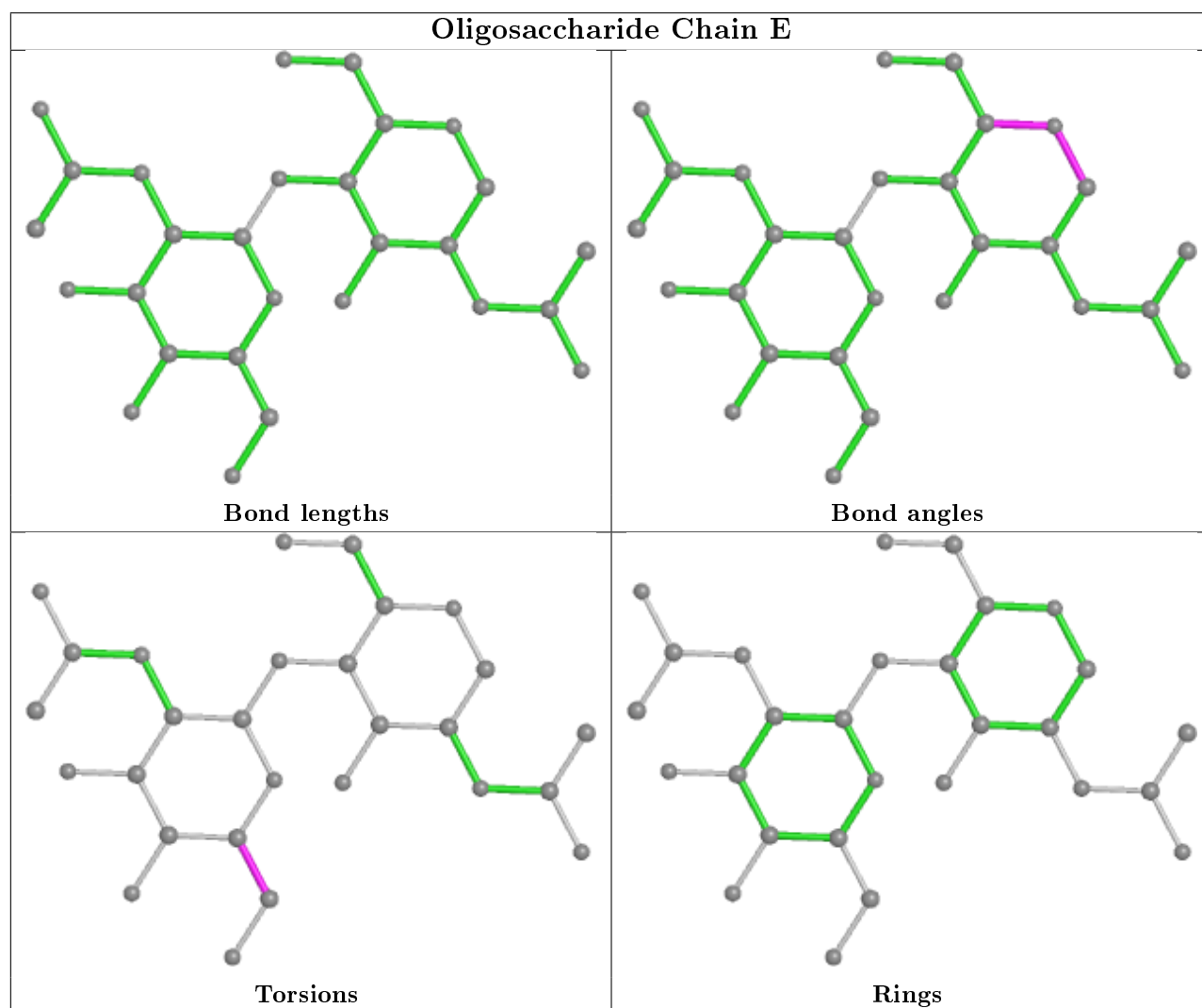
Mol	Chain	Res	Type	Atoms
4	F	2	NAG	O5-C5-C6-O6
4	J	2	NAG	O5-C5-C6-O6
3	H	2	NAG	O5-C5-C6-O6
4	F	2	NAG	C4-C5-C6-O6
4	J	2	NAG	C4-C5-C6-O6

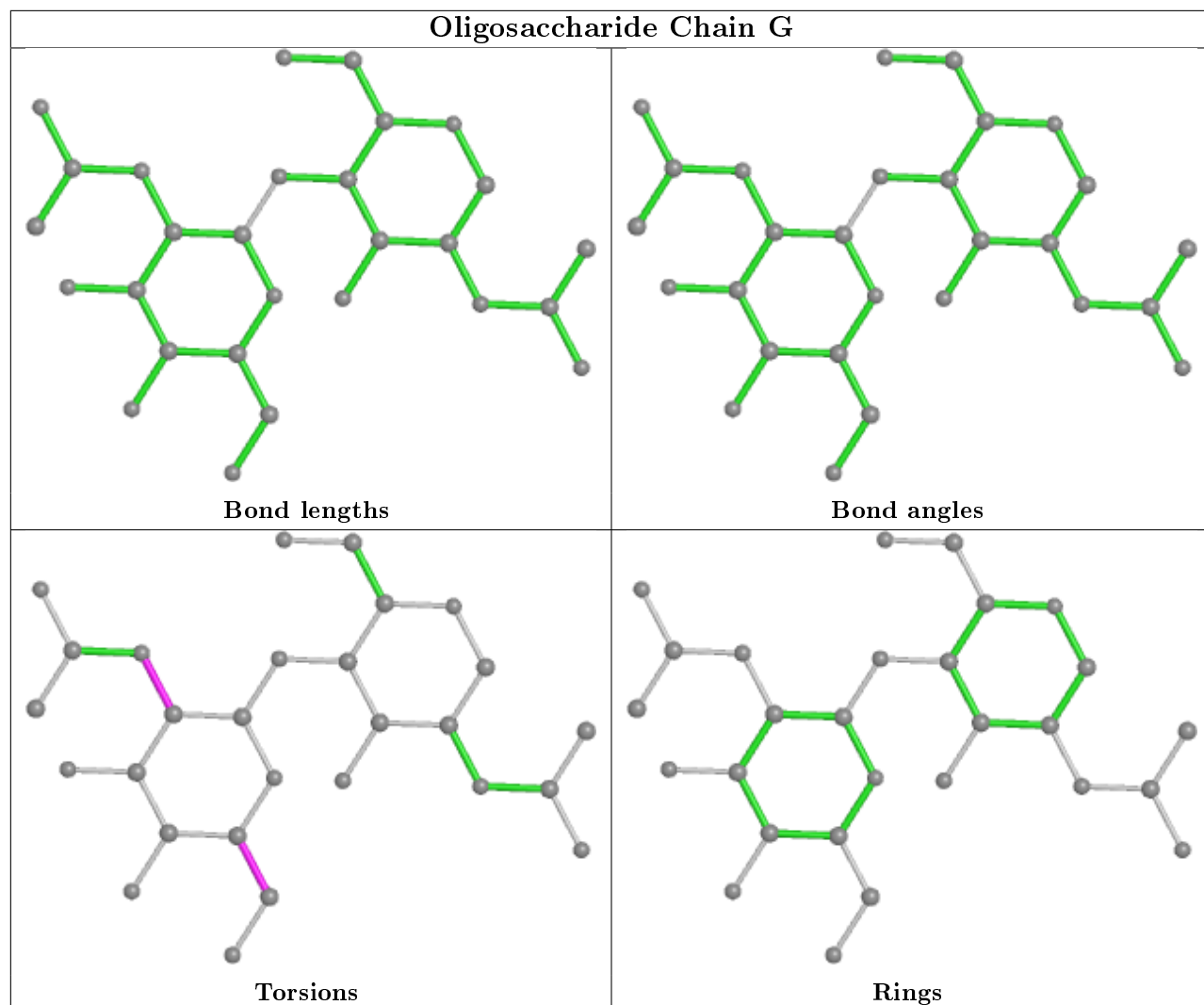
There are no ring outliers.

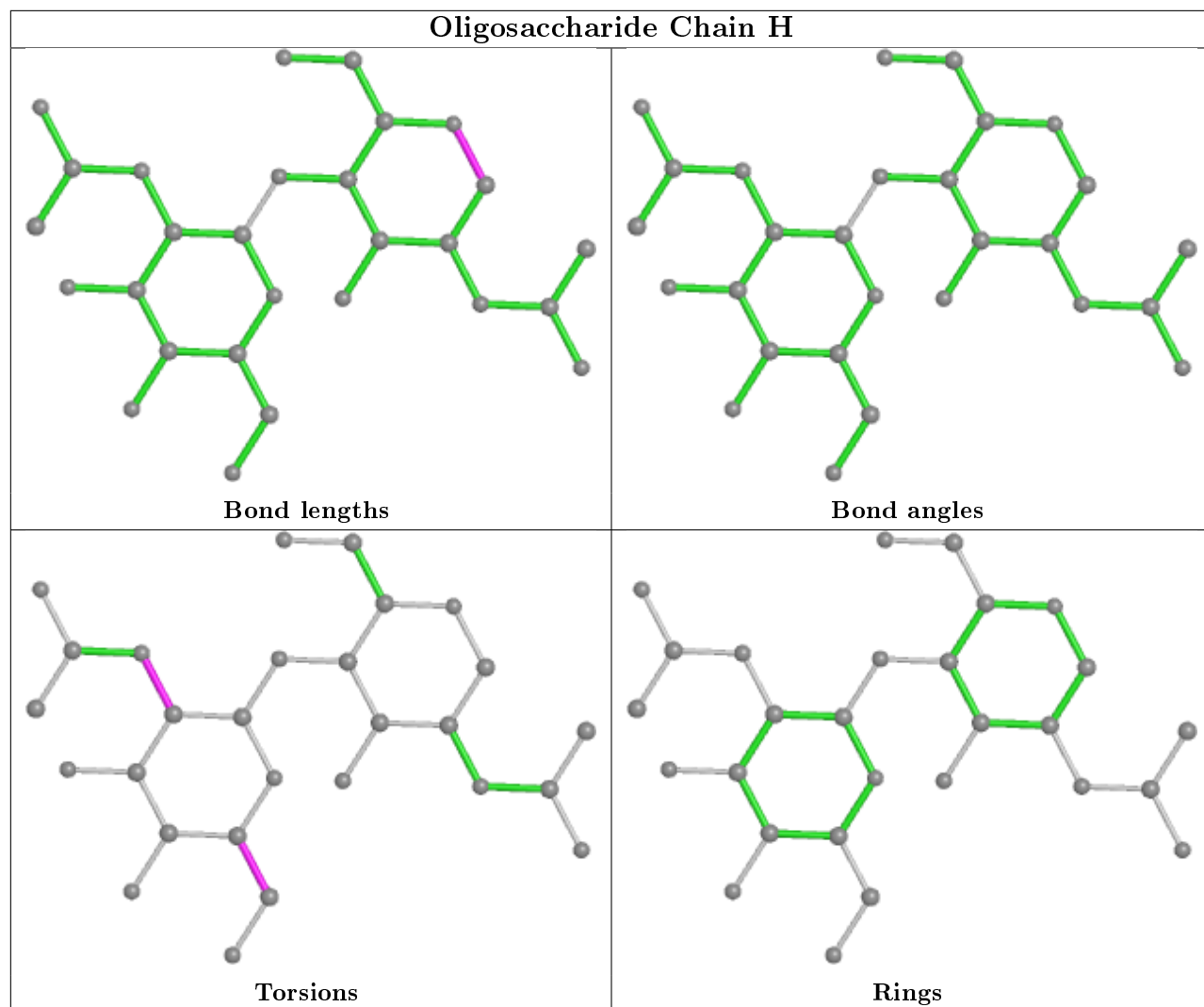
4 monomers are involved in 4 short contacts:

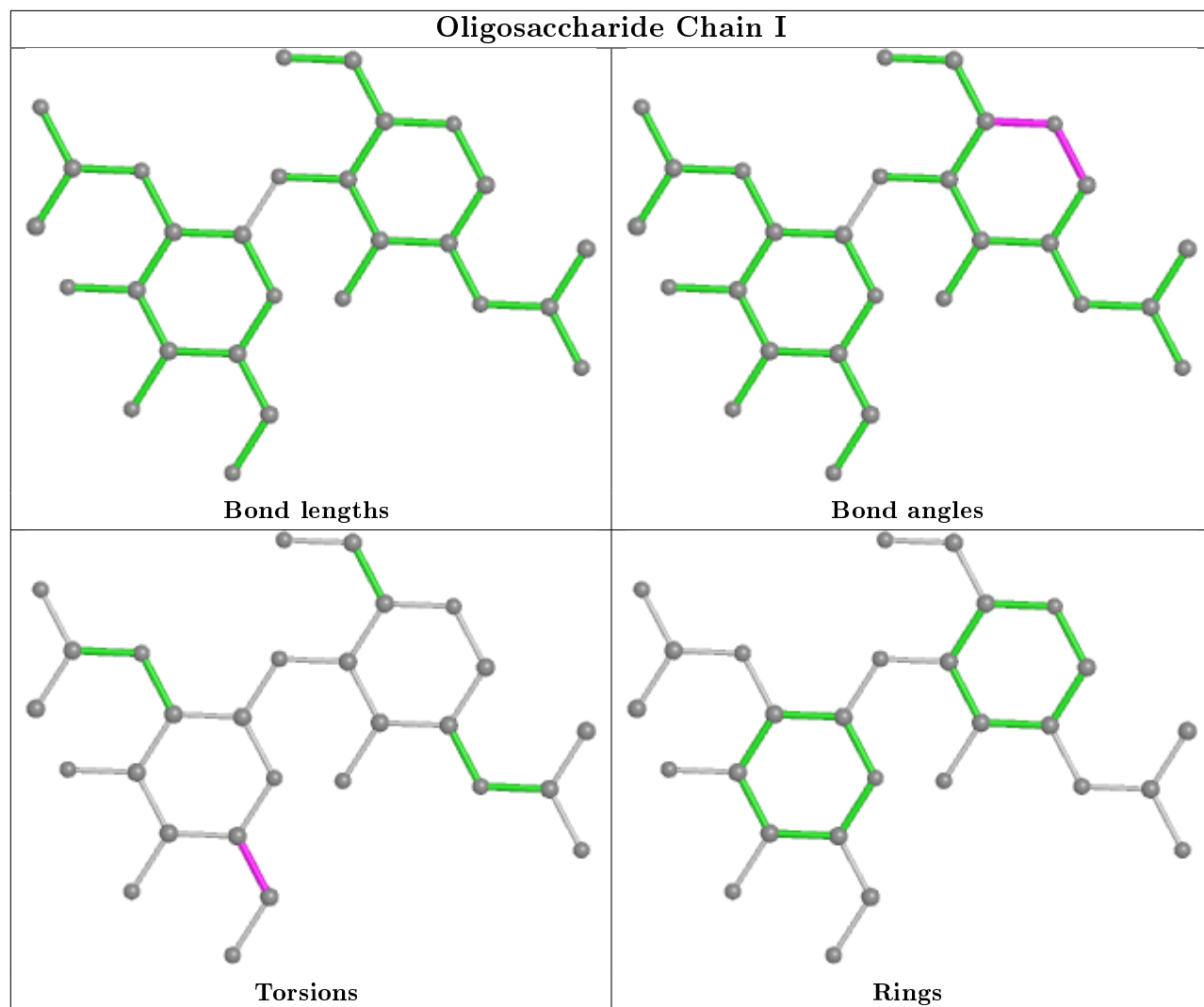
Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	J	2	NAG	2	0
3	G	2	NAG	1	0
3	E	2	NAG	1	0
4	J	3	BMA	1	0

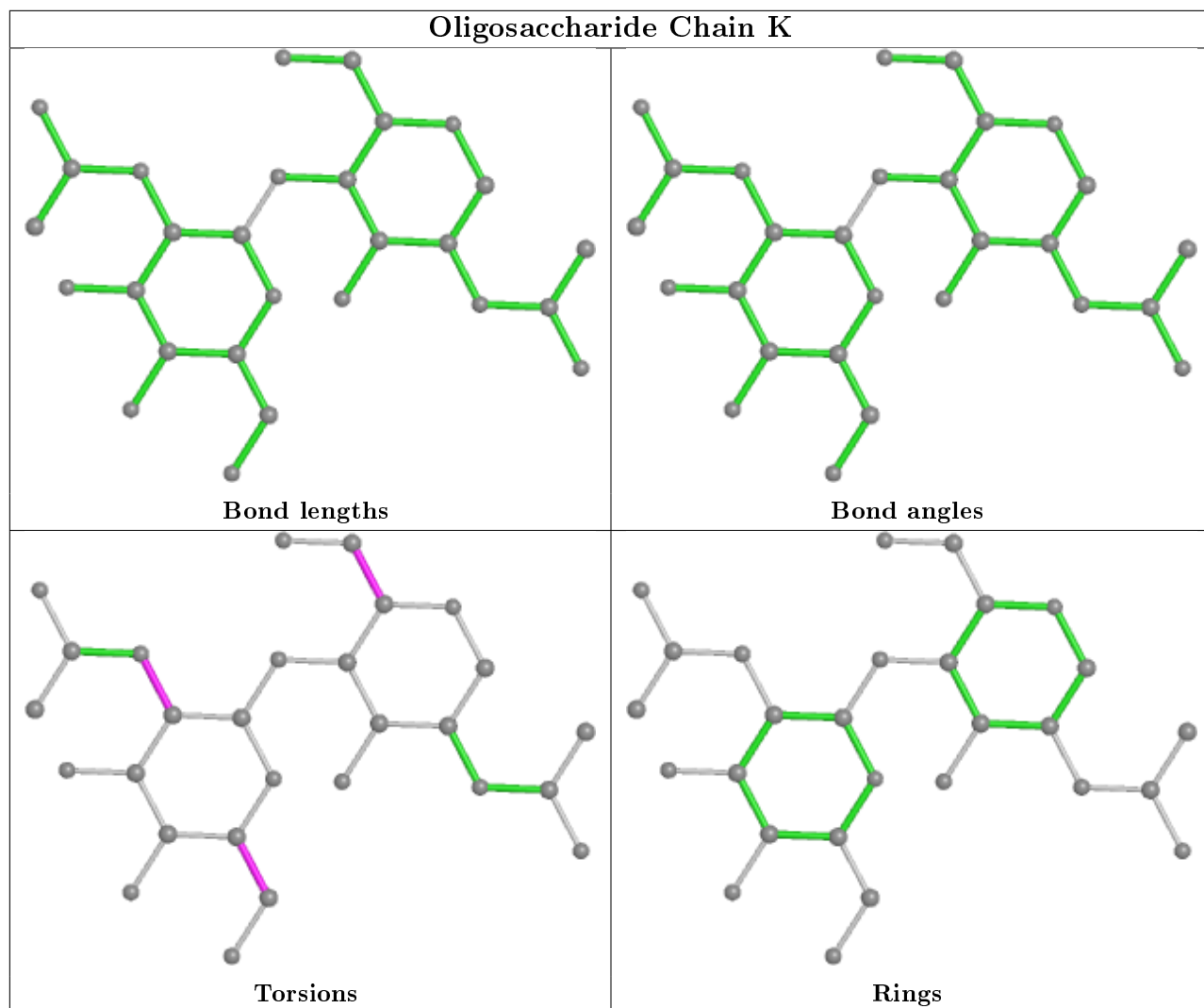
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for oligosaccharide.

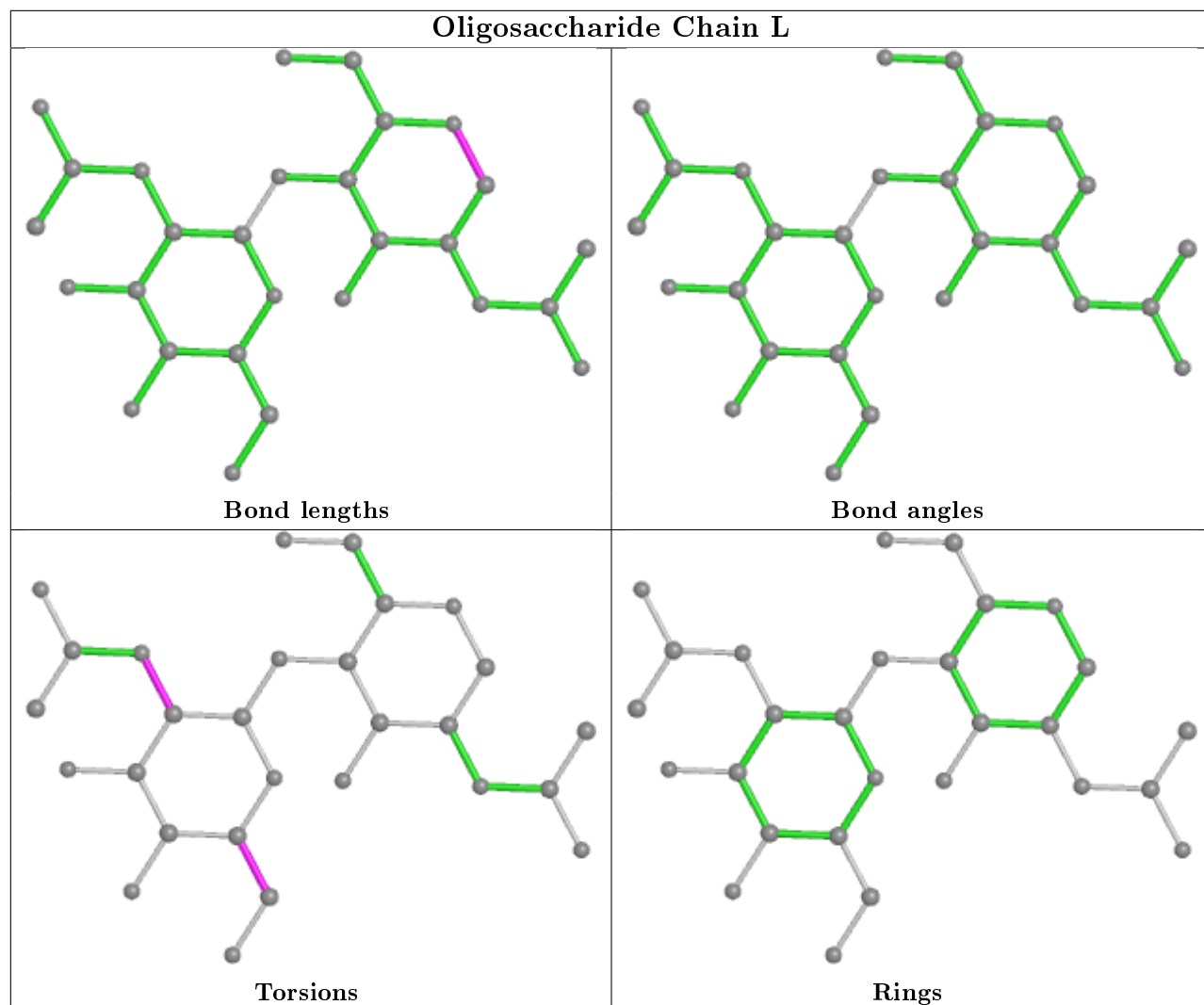


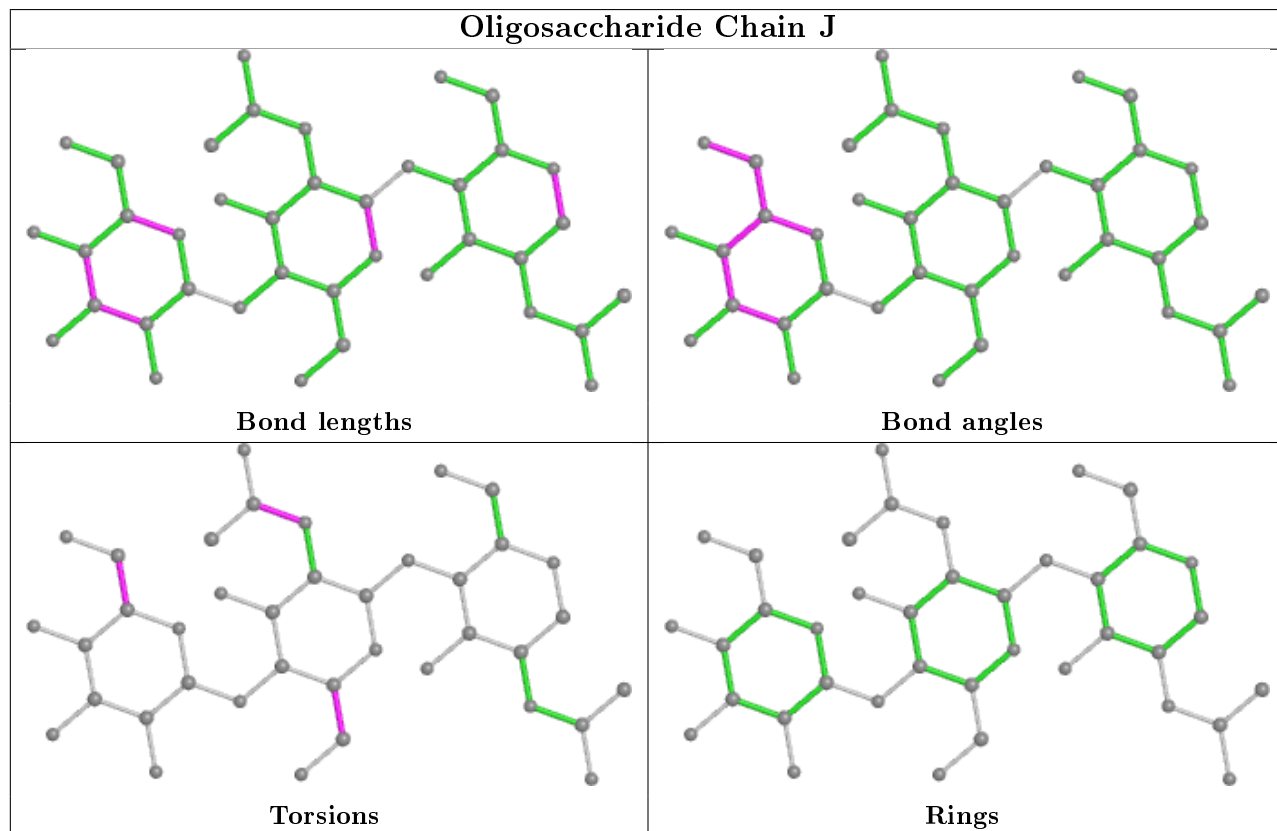
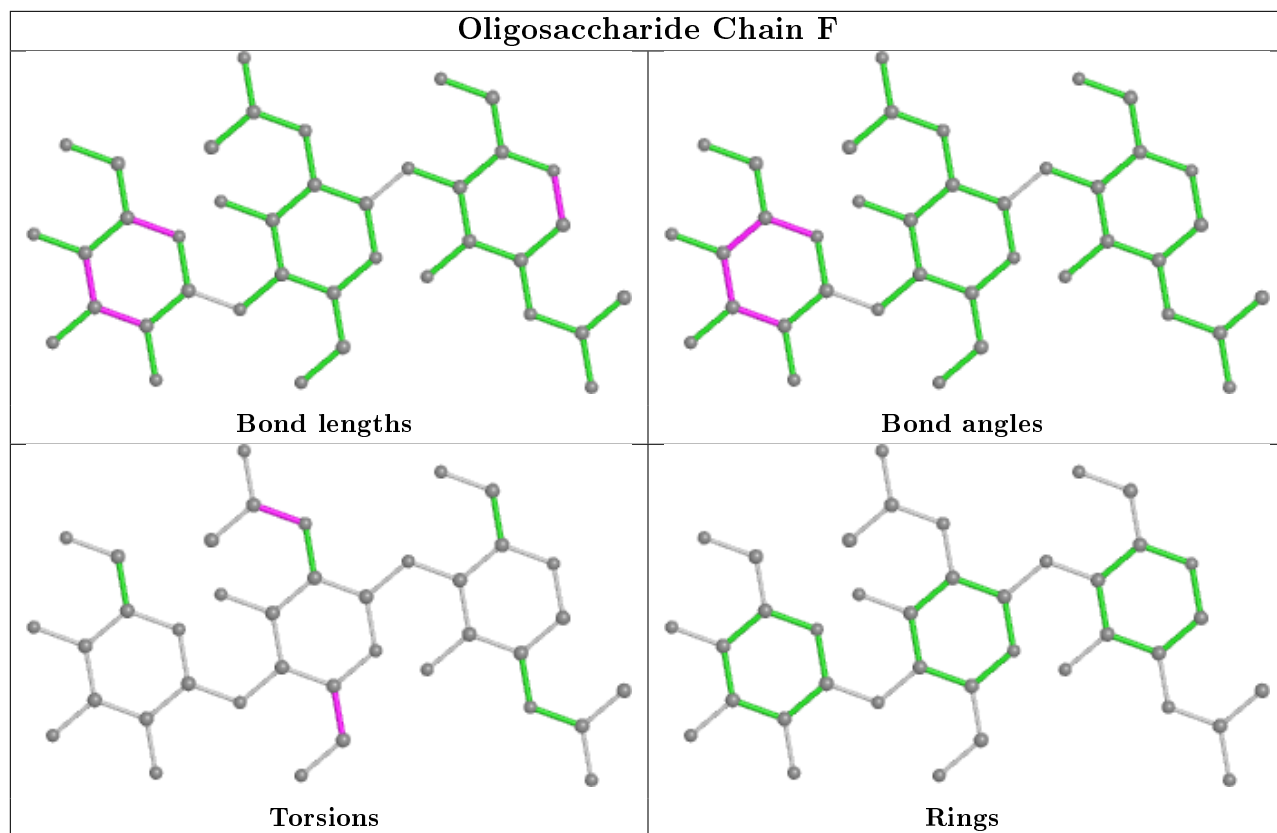












5.6 Ligand geometry

5 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	C	803	1	14,14,15	0.44	0	17,19,21	0.52	0
5	NAG	A	803	1	14,14,15	0.35	0	17,19,21	0.47	0
5	NAG	A	804	1	14,14,15	0.30	0	17,19,21	0.33	0
5	NAG	A	812	1	14,14,15	0.34	0	17,19,21	0.54	0
5	NAG	C	804	1	14,14,15	0.32	0	17,19,21	0.36	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	C	803	1	-	0/6/23/26	0/1/1/1
5	NAG	A	803	1	-	0/6/23/26	0/1/1/1
5	NAG	A	804	1	-	0/6/23/26	0/1/1/1
5	NAG	A	812	1	-	1/6/23/26	0/1/1/1
5	NAG	C	804	1	-	0/6/23/26	0/1/1/1

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (1) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	A	812	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	727/734 (99%)	0.02	12 (1%) 70 66	22, 41, 69, 142	0
1	C	727/734 (99%)	0.04	24 (3%) 46 39	21, 39, 68, 199	0
2	B	208/246 (84%)	0.34	12 (5%) 23 17	34, 55, 96, 119	0
2	D	208/246 (84%)	0.59	18 (8%) 10 7	33, 58, 124, 171	0
All	All	1870/1960 (95%)	0.13	66 (3%) 44 36	21, 43, 84, 199	0

The worst 5 of 66 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	95	PHE	10.3
2	D	213	THR	8.8
2	D	215	THR	8.2
2	D	214	GLY	7.8
1	C	135	TYR	7.2

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

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

6.3 Carbohydrates [i](#)

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

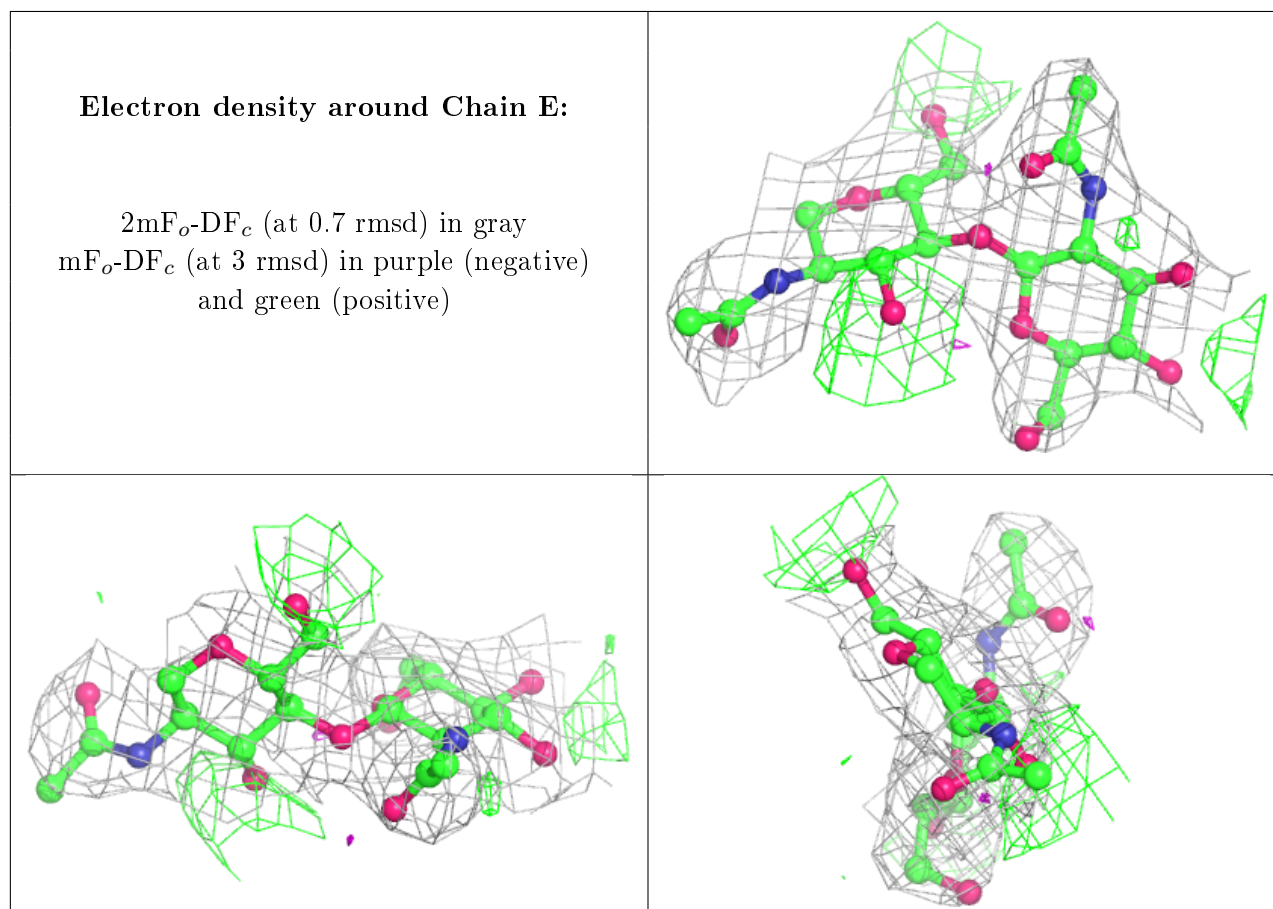
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q < 0.9
4	BMA	J	3	11/12	0.63	0.29	98,109,123,126	0
3	NAG	L	2	14/15	0.77	0.32	76,91,94,95	0

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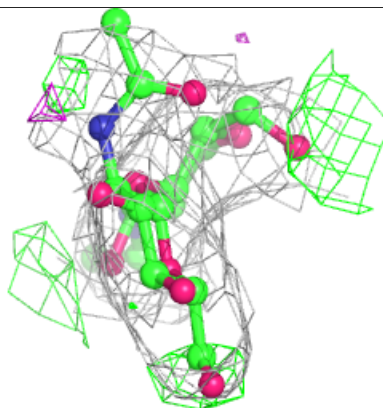
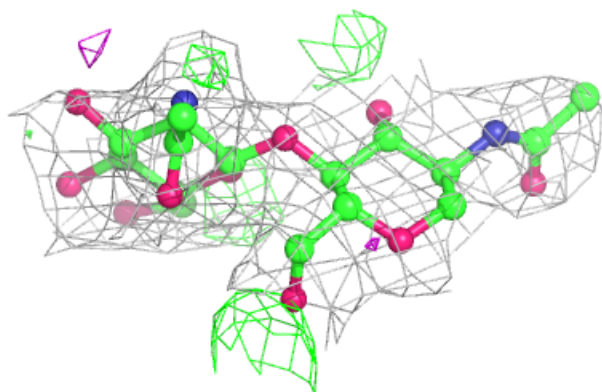
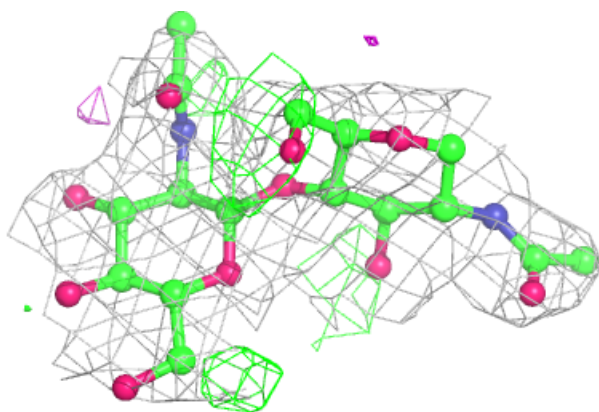
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
3	NAG	E	1	14/15	0.79	0.16	60,63,69,73	0
3	NAG	I	1	14/15	0.83	0.17	65,81,92,102	0
4	BMA	F	3	11/12	0.83	0.31	110,122,127,129	0
3	NAG	K	2	14/15	0.83	0.23	51,70,77,92	0
3	NAG	K	1	14/15	0.84	0.16	51,59,62,62	0
4	NAG	J	2	14/15	0.85	0.26	60,71,79,80	0
3	NAG	I	2	14/15	0.86	0.26	105,114,123,127	0
4	NAG	F	2	14/15	0.86	0.17	51,65,79,84	0
3	NAG	G	2	14/15	0.87	0.26	62,73,79,85	0
3	NAG	H	2	14/15	0.88	0.26	80,86,91,94	0
3	NAG	G	1	14/15	0.89	0.19	46,51,59,61	0
3	NAG	L	1	14/15	0.89	0.26	49,58,66,78	0
3	NAG	E	2	14/15	0.92	0.19	50,60,68,70	0
3	NAG	H	1	14/15	0.94	0.20	57,67,76,82	0
4	NAG	F	1	14/15	0.96	0.16	33,39,54,59	0
4	NAG	J	1	14/15	0.96	0.11	23,40,52,59	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.

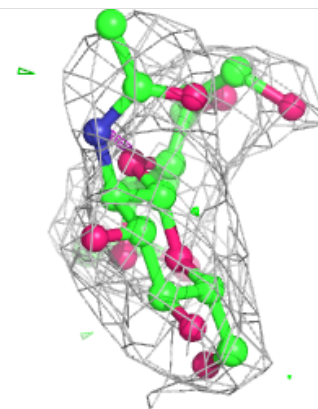
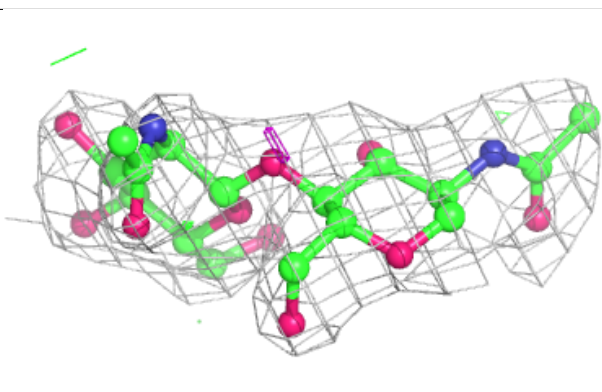
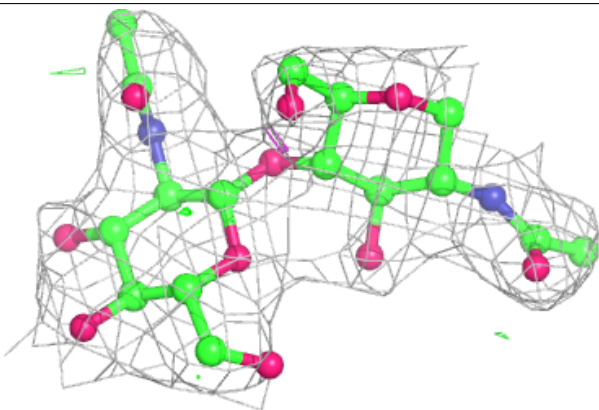


Electron density around Chain G:

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

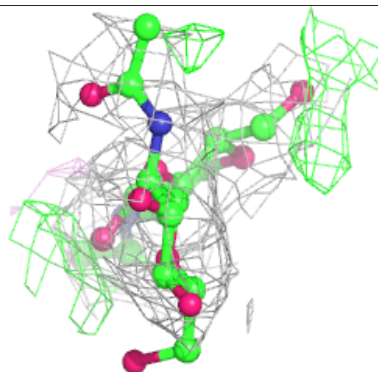
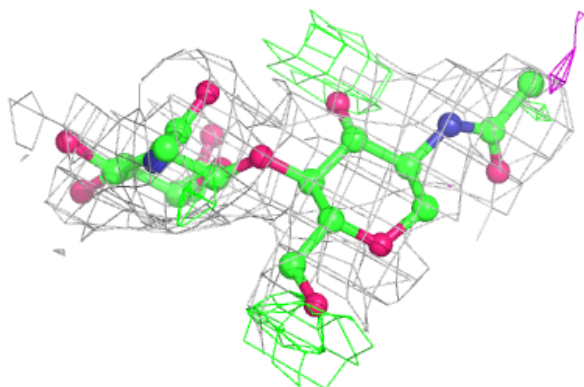
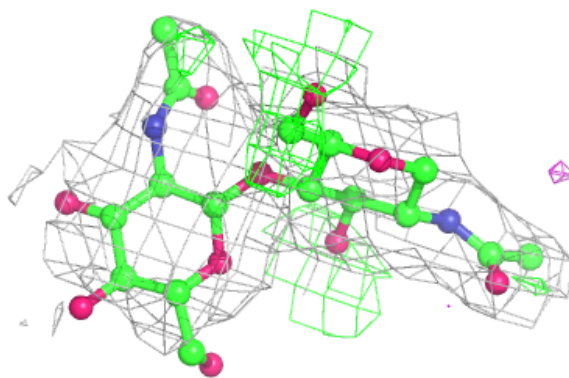
**Electron density around Chain H:**

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

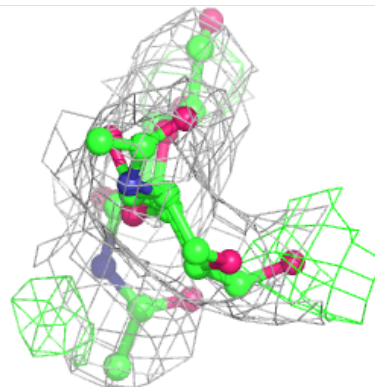
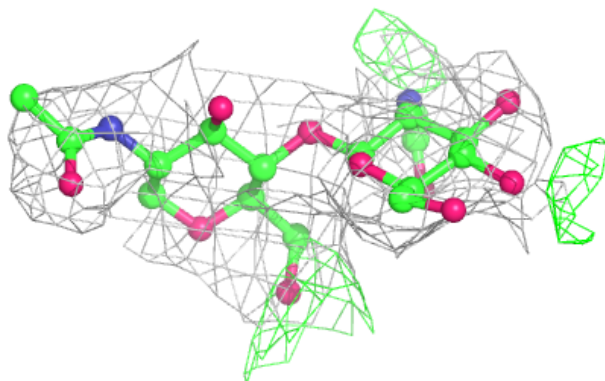
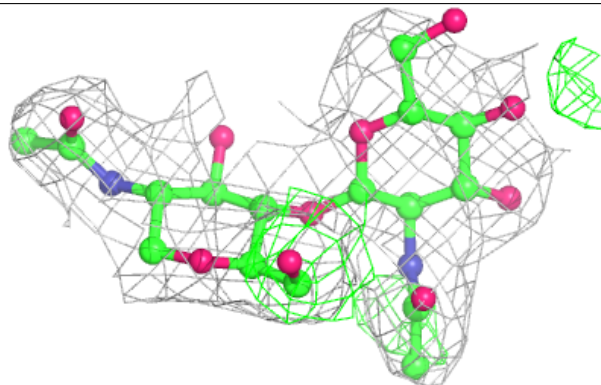


Electron density around Chain I:

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

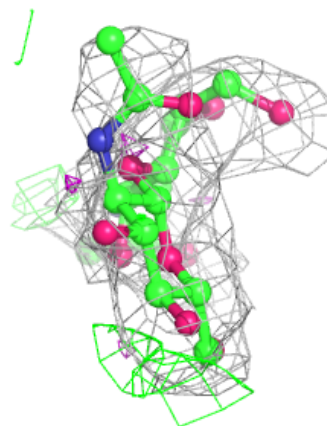
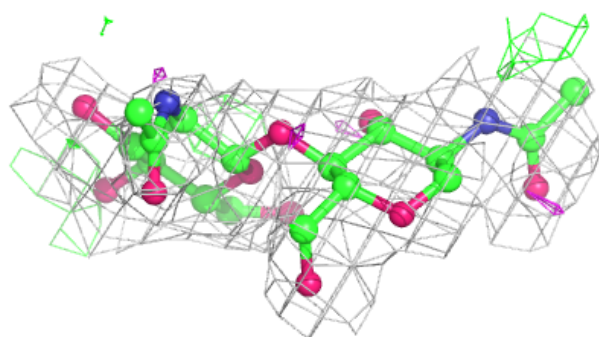
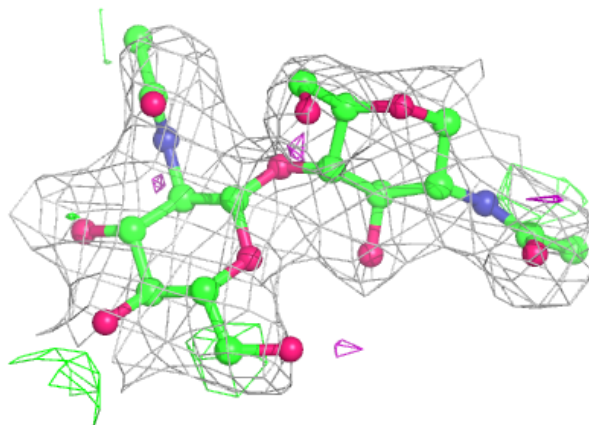
**Electron density around Chain K:**

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

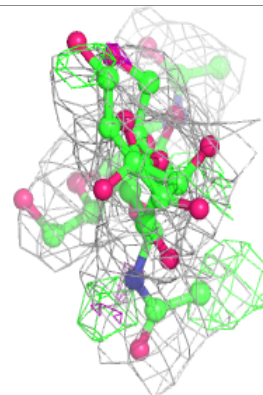
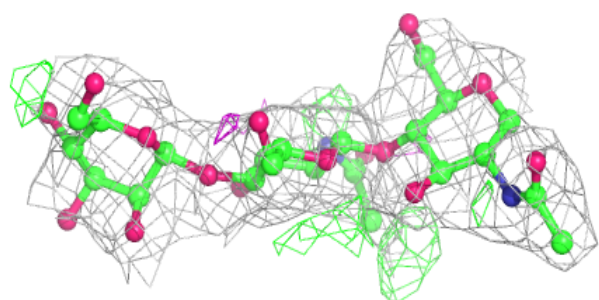
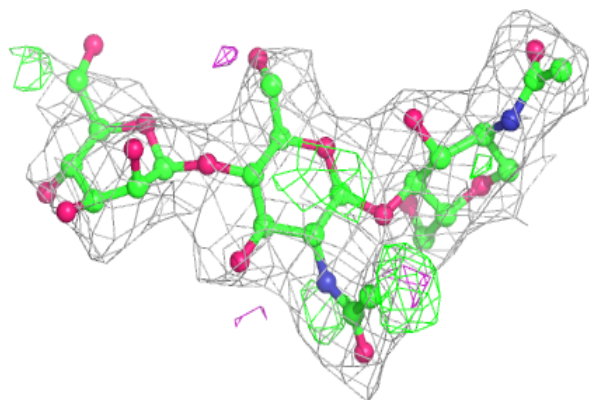


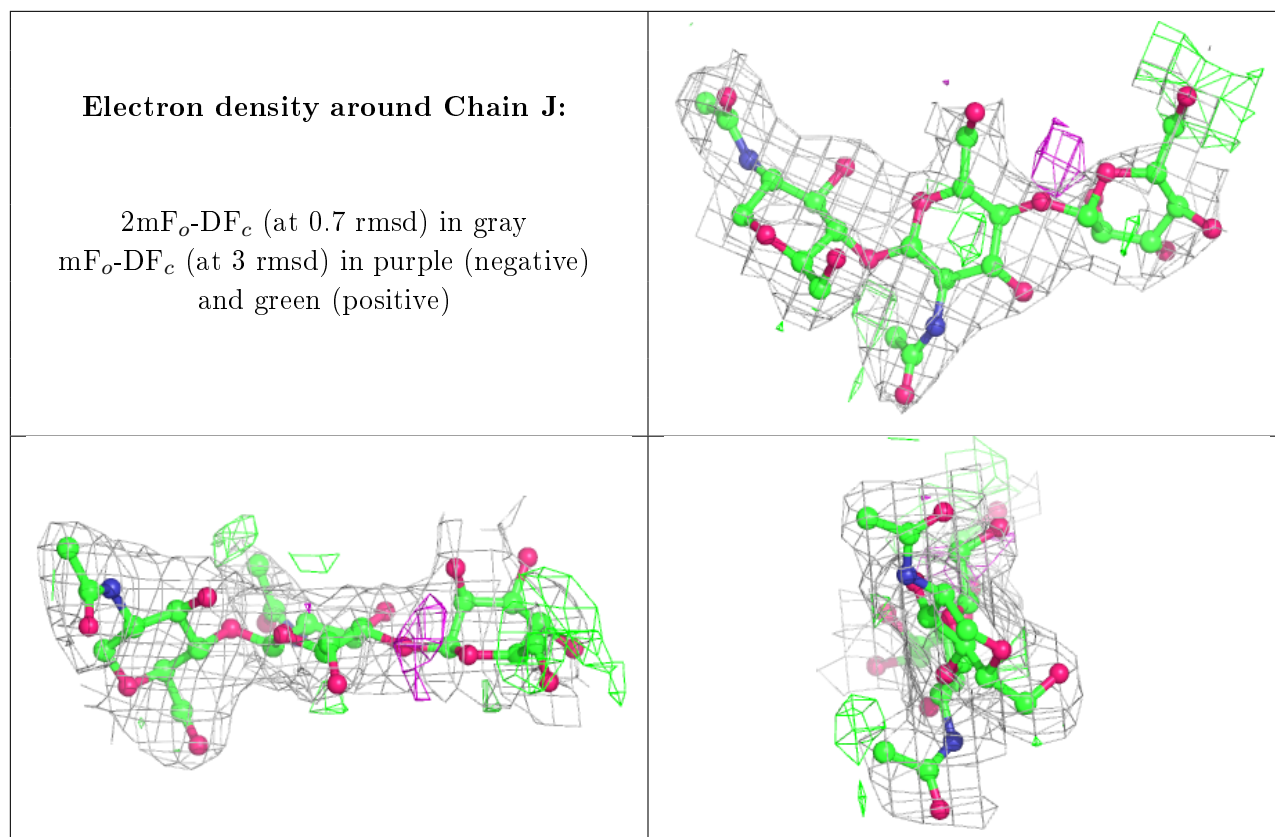
Electron density around Chain L:

$2mF_o-DF_c$ (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)





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	A	812	14/15	0.77	0.28	98,105,114,114	0
5	NAG	A	803	14/15	0.81	0.24	59,82,88,89	0
5	NAG	C	804	14/15	0.83	0.22	71,81,90,92	0
5	NAG	A	804	14/15	0.87	0.15	61,70,78,80	0
5	NAG	C	803	14/15	0.90	0.27	61,80,84,86	0

6.5 Other polymers [i](#)

There are no such residues in this entry.