



Full wwPDB X-ray Structure Validation Report ⓘ

Aug 7, 2020 – 03:17 AM BST

PDB ID : 5E62
Title : HEF-mut with Tr323 complex
Authors : Song, H.; Qi, J.; Shi, Y.; Gao, G.F.
Deposited on : 2015-10-09
Resolution : 2.20 Å(reported)

This is a Full wwPDB X-ray Structure Validation 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

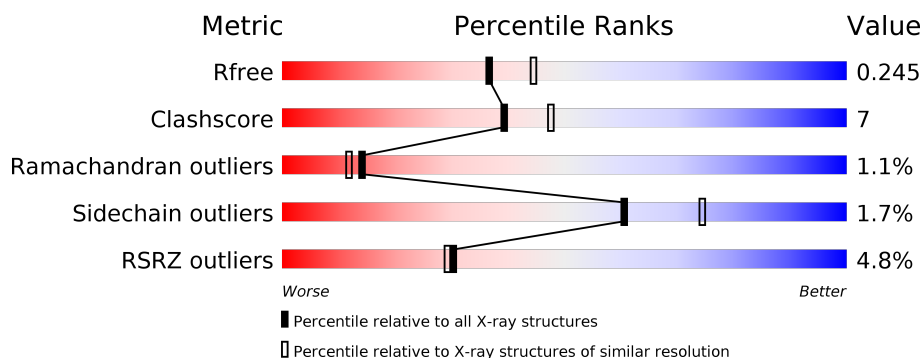
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.20 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	4898 (2.20-2.20)
Clashscore	141614	5594 (2.20-2.20)
Ramachandran outliers	138981	5503 (2.20-2.20)
Sidechain outliers	138945	5504 (2.20-2.20)
RSRZ outliers	127900	4800 (2.20-2.20)

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	427	<div> <div>5%</div> <div>85%</div> <div>13%</div> <div>..</div> </div>
1	C	427	<div> <div>88%</div> <div>11%</div> <div>.</div> </div>
2	B	149	<div> <div>13%</div> <div>75%</div> <div>23%</div> <div>.</div> </div>
2	D	149	<div> <div>9%</div> <div>78%</div> <div>19%</div> <div>.</div> </div>
3	E	6	<div> <div>83%</div> <div>17%</div> </div>
3	F	6	<div> <div>67%</div> <div>33%</div> </div>

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Mol	Chain	Length	Quality of chain
3	J	6	 33% 67%
4	G	2	 50% 50%
4	I	2	 50% 50%
4	K	2	 50% 50%
4	L	2	 50% 50%
4	N	2	 50% 50%
5	H	3	 100%
5	M	3	 33% 67%

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

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
5	Z3Q	H	1	-	-	-	X
5	Z3Q	M	1	-	-	-	X

2 Entry composition [i](#)

There are 7 unique types of molecules in this entry. The entry contains 9821 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-esterase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	424	Total	C	N	O	S	0	0	0
			3240	2039	540	639	22			
1	C	424	Total	C	N	O	S	0	0	0
			3240	2039	540	639	22			

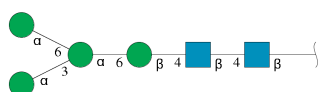
There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	57	ALA	SER	engineered mutation	UNP K9LG83
A	356	ALA	ASP	engineered mutation	UNP K9LG83
A	359	ALA	HIS	engineered mutation	UNP K9LG83
C	57	ALA	SER	engineered mutation	UNP K9LG83
C	356	ALA	ASP	engineered mutation	UNP K9LG83
C	359	ALA	HIS	engineered mutation	UNP K9LG83

- Molecule 2 is a protein called Hemagglutinin-esterase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	149	Total	C	N	O	S	0	0	0
			1110	695	194	217	4			
2	D	149	Total	C	N	O	S	0	0	0
			1110	695	194	217	4			

- Molecule 3 is an oligosaccharide called alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-6)]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	E	6	Total	C	N	O	0	0	0
			72	40	2	30			
3	F	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			

- Molecule 4 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
4	G	2	Total	C	N	O	0	0	0
			28	16	2	10			
4	I	2	Total	C	N	O	0	0	0
			28	16	2	10			
4	K	2	Total	C	N	O	0	0	0
			28	16	2	10			
4	L	2	Total	C	N	O	0	0	0
			28	16	2	10			
4	N	2	Total	C	N	O	0	0	0
			28	16	2	10			

- Molecule 5 is an oligosaccharide called 9-O-acetyl-5-acetamido-3,5-dideoxy-D-glycero-alpha-D-galacto-non-2-ulopyranosonic acid-(2-3)-beta-D-galactopyranose-(1-3)-2-azidoethyl 2-acetamido-2-deoxy-beta-D-glucopyranoside.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
5	H	3	Total	C	N	O	0	0	0
			54	29	5	20			
5	M	3	Total	C	N	O	0	0	0
			54	29	5	20			

- Molecule 6 is 2-acetamido-2-deoxy-beta-D-glucopyranose (three-letter code: NAG) (formula: C₈H₁₅NO₆).



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

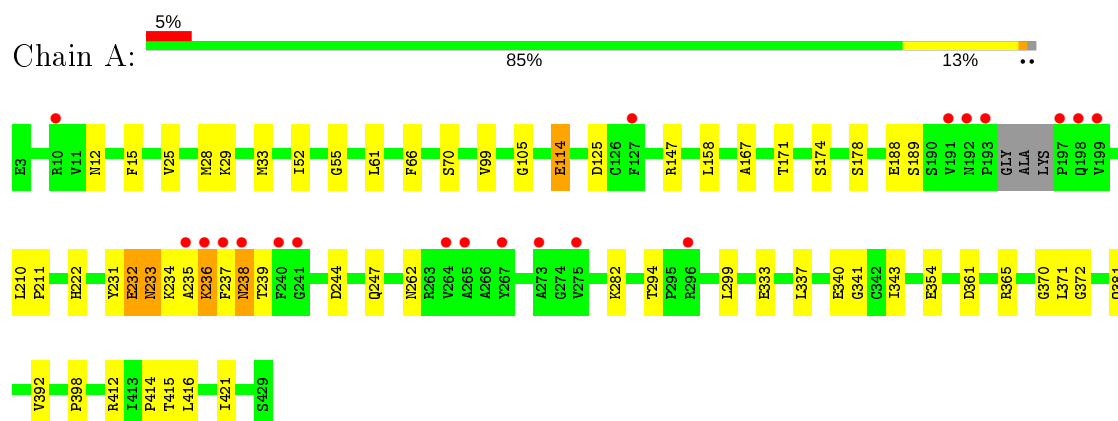
- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	239	Total	O	0	0
			239	239		
7	B	66	Total	O	0	0
			66	66		
7	C	280	Total	O	0	0
			280	280		
7	D	44	Total	O	0	0
			44	44		

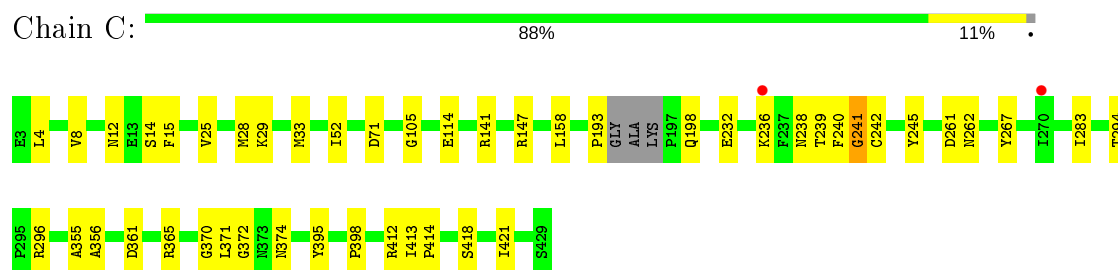
3 Residue-property plots

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.

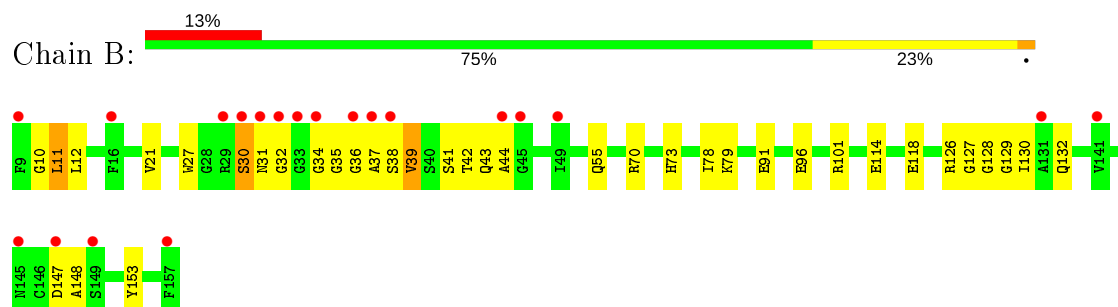
• Molecule 1: Hemagglutinin-esterase



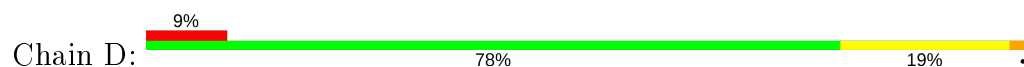
• Molecule 1: Hemagglutinin-esterase



• Molecule 2: Hemagglutinin-esterase



• Molecule 2: Hemagglutinin-esterase





- Molecule 3: alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-6)]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 E: 83% 17%



- Molecule 3: alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-6)]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: 67% 33%



- Molecule 3: alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-6)]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: 33% 67%



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

Chain G: 50% 50%



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

Chain I: 50% 50%



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

Chain K: 50% 50%



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

Chain L:



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

Chain N:



- Molecule 5: 9-O-acetyl-5-acetamido-3,5-dideoxy-D-glycero-alpha-D-galacto-non-2-ulopyranosonic acid-(2-3)-beta-D-galactopyranose-(1-3)-2-azidoethyl 2-acetamido-2-deoxy-beta-D-glucopyranoside

Chain H:



- Molecule 5: 9-O-acetyl-5-acetamido-3,5-dideoxy-D-glycero-alpha-D-galacto-non-2-ulopyranosonic acid-(2-3)-beta-D-galactopyranose-(1-3)-2-azidoethyl 2-acetamido-2-deoxy-beta-D-glucopyranoside

Chain M:



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 3	Depositor
Cell constants a, b, c, α , β , γ	164.82Å 164.82Å 164.82Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	49.70 – 2.20 49.70 – 2.20	Depositor EDS
% Data completeness (in resolution range)	100.0 (49.70-2.20) 96.0 (49.70-2.20)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	0.09	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.32 (at 2.20Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.9_1692)	Depositor
R, R_{free}	0.204 , 0.244 0.208 , 0.245	Depositor DCC
R_{free} test set	3665 reflections (4.87%)	wwPDB-VP
Wilson B-factor (Å ²)	35.5	Xtriage
Anisotropy	0.000	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 39.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.023 for l,-k,h	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	9821	wwPDB-VP
Average B, all atoms (Å ²)	48.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.60% 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, Z3Q, 5N6, GAL, MAN, 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	0.48	0/3315	0.63	1/4494 (0.0%)
1	C	0.46	0/3315	0.61	0/4494
2	B	0.44	0/1127	0.62	0/1515
2	D	0.38	0/1127	0.55	0/1515
All	All	0.46	0/8884	0.61	1/12018 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	233	ASN	C-N-CA	-5.47	108.01	121.70

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	3240	0	3114	47	0
1	C	3240	0	3114	38	0
2	B	1110	0	1067	25	0
2	D	1110	0	1067	28	0
3	E	72	0	61	1	0
3	F	72	0	61	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	J	72	0	61	0	0
4	G	28	0	25	1	0
4	I	28	0	25	2	0
4	K	28	0	25	0	0
4	L	28	0	25	1	0
4	N	28	0	25	1	0
5	H	54	0	9	6	0
5	M	54	0	9	2	0
6	B	14	0	13	0	0
6	D	14	0	13	0	0
7	A	239	0	0	5	1
7	B	66	0	0	5	1
7	C	280	0	0	7	1
7	D	44	0	0	7	0
All	All	9821	0	8714	132	2

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 7.

All (132) 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:234:LYS:O	1:A:238:ASN:ND2	1.89	1.06
1:A:244:ASP:OD1	1:A:262:ASN:HB2	1.70	0.92
1:A:29:LYS:NZ	7:A:804:HOH:O	2.05	0.90
5:H:3:5N6:O7	5:H:3:5N6:O10	1.90	0.90
2:D:80:ASN:OD1	7:D:801:HOH:O	1.91	0.87
1:C:193:PRO:O	7:C:801:HOH:O	1.96	0.82
1:A:340:GLU:OE1	7:A:802:HOH:O	1.98	0.81
1:A:114:GLU:OE2	7:A:803:HOH:O	1.99	0.79
2:D:80:ASN:ND2	7:D:802:HOH:O	2.14	0.77
1:C:198:GLN:OE1	7:C:802:HOH:O	2.03	0.77
2:D:30:SER:OG	2:D:31:ASN:ND2	2.19	0.76
1:A:415:THR:O	2:B:101:ARG:NH1	2.19	0.74
7:B:801:HOH:O	4:I:2:NAG:O4	2.06	0.72
1:A:233:ASN:O	1:A:234:LYS:C	2.19	0.72
1:C:240:PHE:O	1:C:296:ARG:NH1	2.23	0.72
5:M:2:GAL:H4	5:M:3:5N6:C1	2.19	0.72
1:C:232:GLU:OE1	1:C:236:LYS:NZ	2.19	0.70
3:F:2:NAG:O3	3:F:3:BMA:H2	1.91	0.70
1:C:261:ASP:OD2	7:C:803:HOH:O	2.08	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:114:GLU:HG2	7:A:803:HOH:O	1.92	0.70
1:C:245:TYR:O	1:C:262:ASN:HB2	1.92	0.70
1:A:232:GLU:HG3	1:A:236:LYS:HG3	1.72	0.69
1:A:236:LYS:O	1:A:239:THR:OG1	2.06	0.69
1:C:238:ASN:HA	1:C:241:GLY:O	1.92	0.69
2:D:30:SER:OG	2:D:31:ASN:N	2.24	0.66
1:A:237:PHE:HZ	1:A:294:THR:HG21	1.61	0.66
3:E:2:NAG:O3	3:E:3:BMA:C1	2.43	0.66
1:C:236:LYS:O	1:C:239:THR:OG1	2.13	0.65
1:C:25:VAL:HG13	1:C:421:ILE:HG23	1.77	0.65
1:A:33:MET:HE1	2:B:101:ARG:HB2	1.79	0.65
1:A:234:LYS:HG3	1:A:238:ASN:HD21	1.61	0.65
2:B:38:SER:OG	2:B:39:VAL:N	2.30	0.63
1:C:412:ARG:NE	7:C:809:HOH:O	2.31	0.63
2:B:101:ARG:NH1	7:B:804:HOH:O	2.31	0.63
1:A:33:MET:HG3	1:A:416:LEU:HG	1.80	0.63
1:C:52:ILE:HD13	1:C:105:GLY:HA3	1.81	0.63
1:A:114:GLU:OE1	1:A:114:GLU:N	2.32	0.62
1:A:361:ASP:O	1:A:365:ARG:HG3	2.00	0.62
2:D:101:ARG:NE	7:D:804:HOH:O	2.25	0.60
1:C:242:CYS:HB2	7:C:901:HOH:O	2.02	0.60
2:D:79:LYS:NZ	7:D:805:HOH:O	2.35	0.60
1:A:233:ASN:O	1:A:235:ALA:N	2.35	0.59
1:C:33:MET:HE2	1:C:414:PRO:HB2	1.84	0.58
1:A:232:GLU:HG3	1:A:236:LYS:CG	2.34	0.58
1:C:418:SER:HB2	2:D:57:ARG:HH11	1.68	0.58
2:B:96:GLU:OE2	7:B:803:HOH:O	2.17	0.58
1:C:8:VAL:HG13	2:D:26:PHE:HB2	1.86	0.56
1:A:238:ASN:H	1:A:238:ASN:HD22	1.52	0.56
2:B:27:TRP:CD1	2:B:44:ALA:HB2	2.43	0.54
1:A:28:MET:SD	4:G:1:NAG:H5	2.48	0.53
2:D:26:PHE:HD1	2:D:43:GLN:H	1.57	0.53
1:A:25:VAL:HG13	1:A:421:ILE:HG23	1.90	0.53
2:B:126:ARG:NE	2:B:153:TYR:O	2.28	0.53
2:B:42:THR:O	2:B:43:GLN:HB2	2.09	0.53
1:C:361:ASP:O	1:C:365:ARG:HG3	2.09	0.52
1:C:147:ARG:NH1	2:D:75:GLU:OE1	2.42	0.52
1:A:333:GLU:OE1	7:A:805:HOH:O	2.19	0.52
1:A:412:ARG:NH2	2:B:91:GLU:HA	2.25	0.51
1:C:12:ASN:HB3	1:C:14:SER:H	1.76	0.51
1:C:294:THR:OG1	7:C:805:HOH:O	2.19	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:4:LEU:H	2:D:30:SER:CB	2.24	0.51
1:A:354:GLU:O	1:C:29:LYS:HE2	2.10	0.51
1:A:61:LEU:HA	1:A:66:PHE:CD1	2.47	0.50
1:A:371:LEU:N	1:A:372:GLY:HA3	2.27	0.50
2:D:27:TRP:CD1	2:D:44:ALA:HB2	2.47	0.50
1:C:4:LEU:HB3	2:D:30:SER:HB2	1.93	0.50
1:A:167:ALA:HA	1:A:188:GLU:HG3	1.93	0.49
1:C:232:GLU:HB3	1:C:236:LYS:HE2	1.94	0.49
1:A:55:GLY:HA2	1:A:114:GLU:HB3	1.94	0.49
1:C:371:LEU:N	1:C:372:GLY:HA3	2.28	0.49
5:H:2:GAL:H4	5:H:3:5N6:C1	2.43	0.48
1:A:52:ILE:HD13	1:A:105:GLY:HA3	1.95	0.48
1:C:240:PHE:H	1:C:241:GLY:HA3	1.78	0.48
1:A:12:ASN:H	1:A:15:PHE:HD1	1.60	0.47
5:H:2:GAL:C4	5:H:3:5N6:C1	2.93	0.47
1:A:147:ARG:HD3	1:A:341:GLY:HA3	1.96	0.47
2:B:10:GLY:O	2:B:12:LEU:N	2.40	0.47
2:B:37:ALA:O	2:B:38:SER:HB3	2.15	0.47
2:D:101:ARG:NH2	7:D:804:HOH:O	2.47	0.46
2:D:34:GLY:HA2	2:D:35:GLY:HA2	1.68	0.46
5:M:2:GAL:C4	5:M:3:5N6:C1	2.86	0.46
7:D:803:HOH:O	4:N:2:NAG:O4	2.20	0.46
2:B:34:GLY:HA2	2:B:35:GLY:HA2	1.67	0.46
2:D:149:SER:O	2:D:153:TYR:N	2.49	0.46
2:D:147:ASP:OD2	2:D:148:ALA:N	2.40	0.45
2:D:51:LYS:O	2:D:51:LYS:HD2	2.16	0.45
1:C:33:MET:CE	2:D:101:ARG:HB2	2.46	0.45
1:C:52:ILE:CD1	1:C:105:GLY:HA3	2.47	0.45
1:C:370:GLY:C	1:C:372:GLY:HA3	2.36	0.45
1:A:231:TYR:OH	5:H:3:5N6:OBJ	2.28	0.45
1:A:343:ILE:HG23	1:A:381:GLN:HA	1.99	0.45
1:A:412:ARG:HH22	2:B:91:GLU:HG3	1.82	0.45
1:C:71:ASP:HA	1:C:365:ARG:HG2	1.98	0.45
2:B:73:HIS:HE1	7:B:806:HOH:O	2.00	0.45
1:A:282:LYS:HB2	1:A:282:LYS:HE2	1.86	0.44
2:B:114:GLU:O	2:B:118:GLU:HG3	2.17	0.44
2:B:31:ASN:HA	2:B:36:GLY:HA3	1.99	0.44
1:C:395:TYR:HD2	2:D:86:GLU:HG2	1.82	0.44
1:A:412:ARG:HH22	2:B:91:GLU:HA	1.83	0.44
1:C:413:ILE:HA	1:C:414:PRO:HD3	1.87	0.44
1:A:171:THR:HG22	1:A:178:SER:HA	1.99	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:12:ASN:H	1:C:15:PHE:HD1	1.65	0.43
2:D:128:GLY:HA2	2:D:129:GLY:HA2	1.52	0.43
2:B:30:SER:HB3	2:B:31:ASN:H	1.50	0.43
2:B:55:GLN:NE2	4:I:2:NAG:H83	2.33	0.43
1:A:232:GLU:HG2	1:A:232:GLU:H	1.50	0.42
1:A:337:LEU:O	2:B:79:LYS:HD3	2.19	0.42
1:A:125:ASP:HB3	1:A:174:SER:O	2.19	0.42
1:A:211:PRO:HD2	1:A:222:HIS:CE1	2.54	0.42
2:B:128:GLY:HA2	2:B:129:GLY:HA2	1.71	0.42
2:B:127:GLY:O	2:B:130:ILE:N	2.52	0.42
2:B:70:ARG:NH1	7:B:807:HOH:O	2.52	0.42
1:C:262:ASN:ND2	1:C:267:TYR:CD1	2.88	0.42
1:C:374:ASN:ND2	7:C:826:HOH:O	2.53	0.42
5:H:1:Z3Q:C8	5:H:2:GAL:H2	2.50	0.41
1:C:395:TYR:CD2	2:D:86:GLU:HG2	2.55	0.41
2:B:147:ASP:OD2	2:B:148:ALA:N	2.44	0.41
2:D:38:SER:OG	2:D:39:VAL:N	2.53	0.41
1:A:210:LEU:HB3	1:A:222:HIS:CD2	2.56	0.41
2:D:142:ASP:C	2:D:144:GLU:H	2.24	0.41
1:A:370:GLY:C	1:A:372:GLY:HA3	2.40	0.41
1:A:247:GLN:NE2	1:A:299:LEU:O	2.51	0.41
1:C:398:PRO:HB2	2:D:78:ILE:O	2.21	0.41
1:A:70:SER:O	1:A:365:ARG:HD3	2.21	0.41
1:C:141:ARG:HD2	1:C:283:ILE:HG13	2.01	0.41
1:C:28:MET:SD	4:L:1:NAG:H5	2.61	0.41
2:D:126:ARG:HD3	2:D:154:ILE:HA	2.03	0.41
2:D:42:THR:O	2:D:42:THR:HG23	2.21	0.40
2:D:51:LYS:HG3	7:D:827:HOH:O	2.21	0.40
1:A:33:MET:CE	1:A:414:PRO:HB2	2.50	0.40
1:A:398:PRO:HB2	2:B:78:ILE:O	2.21	0.40
5:H:3:5N6:O10	5:H:3:5N6:C7	2.70	0.40

All (2) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
7:A:977:HOH:O	7:C:1012:HOH:O[3_554]	1.99	0.21
7:B:807:HOH:O	7:B:829:HOH:O[10_554]	2.19	0.01

5.3 Torsion angles

5.3.1 Protein backbone

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	420/427 (98%)	399 (95%)	21 (5%)	0	100	100
1	C	420/427 (98%)	402 (96%)	15 (4%)	3 (1%)	22	22
2	B	147/149 (99%)	124 (84%)	17 (12%)	6 (4%)	3	1
2	D	147/149 (99%)	127 (86%)	16 (11%)	4 (3%)	5	2
All	All	1134/1152 (98%)	1052 (93%)	69 (6%)	13 (1%)	14	12

All (13) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	11	LEU
1	C	355	ALA
2	D	128	GLY
2	B	21	VAL
2	B	41	SER
2	B	132	GLN
2	D	41	SER
2	D	132	GLN
2	D	43	GLN
1	C	356	ALA
2	B	32	GLY
1	C	241	GLY
2	B	39	VAL

5.3.2 Protein sidechains

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	357/358 (100%)	349 (98%)	8 (2%)	52	65
1	C	357/358 (100%)	355 (99%)	2 (1%)	86	93
2	B	109/109 (100%)	107 (98%)	2 (2%)	59	72
2	D	109/109 (100%)	105 (96%)	4 (4%)	34	43
All	All	932/934 (100%)	916 (98%)	16 (2%)	60	74

All (16) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	99	VAL
1	A	114	GLU
1	A	158	LEU
1	A	189	SER
1	A	232	GLU
1	A	236	LYS
1	A	238	ASN
1	A	392	VAL
2	B	11	LEU
2	B	30	SER
1	C	114	GLU
1	C	158	LEU
2	D	9	PHE
2	D	11	LEU
2	D	126	ARG
2	D	144	GLU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	238	ASN
2	B	156	ASN
2	D	31	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

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

5.5 Carbohydrates ⓘ

34 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	0.42	0	17,19,21	0.87	2 (11%)
3	NAG	E	2	3	14,14,15	0.30	0	17,19,21	0.69	0
3	BMA	E	3	3	11,11,12	0.92	1 (9%)	15,15,17	2.84	5 (33%)
3	MAN	E	4	3	11,11,12	0.39	0	15,15,17	1.28	1 (6%)
3	MAN	E	5	3	11,11,12	0.34	0	15,15,17	1.43	2 (13%)
3	MAN	E	6	3	11,11,12	0.39	0	15,15,17	1.19	1 (6%)
3	NAG	F	1	1,3	14,14,15	0.43	0	17,19,21	0.95	1 (5%)
3	NAG	F	2	3	14,14,15	0.55	0	17,19,21	1.67	4 (23%)
3	BMA	F	3	3	11,11,12	0.60	0	15,15,17	3.23	6 (40%)
3	MAN	F	4	3	11,11,12	0.49	0	15,15,17	2.07	3 (20%)
3	MAN	F	5	3	11,11,12	0.34	0	15,15,17	0.90	1 (6%)
3	MAN	F	6	3	11,11,12	0.49	0	15,15,17	1.13	1 (6%)
4	NAG	G	1	1,4	14,14,15	0.77	1 (7%)	17,19,21	1.35	2 (11%)
4	NAG	G	2	4	14,14,15	0.30	0	17,19,21	1.40	3 (17%)
5	Z3Q	H	1	5	19,20,20	1.72	2 (10%)	24,26,26	3.83	7 (29%)
5	GAL	H	2	5	11,11,12	1.02	1 (9%)	15,15,17	2.27	4 (26%)
5	5N6	H	3	5	20,23,24	1.92	5 (25%)	25,32,35	1.70	8 (32%)
4	NAG	I	1	2,4	14,14,15	0.44	0	17,19,21	0.90	0
4	NAG	I	2	4	14,14,15	0.39	0	17,19,21	1.36	2 (11%)
3	NAG	J	1	1,3	14,14,15	0.45	0	17,19,21	0.79	0
3	NAG	J	2	3	14,14,15	0.50	0	17,19,21	1.23	1 (5%)
3	BMA	J	3	3	11,11,12	0.49	0	15,15,17	1.03	1 (6%)
3	MAN	J	4	3	11,11,12	0.48	0	15,15,17	1.66	4 (26%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	MAN	J	5	3	11,11,12	0.52	0	15,15,17	1.12	1 (6%)
3	MAN	J	6	3	11,11,12	0.39	0	15,15,17	0.73	0
4	NAG	K	1	1,4	14,14,15	0.50	0	17,19,21	1.06	2 (11%)
4	NAG	K	2	4	14,14,15	0.50	0	17,19,21	0.84	0
4	NAG	L	1	1,4	14,14,15	0.38	0	17,19,21	1.09	1 (5%)
4	NAG	L	2	4	14,14,15	0.36	0	17,19,21	1.31	1 (5%)
5	Z3Q	M	1	5	19,20,20	1.20	2 (10%)	24,26,26	1.06	2 (8%)
5	GAL	M	2	5	11,11,12	1.19	0	15,15,17	1.33	3 (20%)
5	5N6	M	3	5	20,23,24	2.52	9 (45%)	25,32,35	1.11	1 (4%)
4	NAG	N	1	2,4	14,14,15	0.45	0	17,19,21	0.93	2 (11%)
4	NAG	N	2	4	14,14,15	0.48	0	17,19,21	1.41	2 (11%)

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	-	0/6/23/26	0/1/1/1
3	BMA	E	3	3	-	0/2/19/22	0/1/1/1
3	MAN	E	4	3	-	0/2/19/22	0/1/1/1
3	MAN	E	5	3	-	2/2/19/22	0/1/1/1
3	MAN	E	6	3	-	0/2/19/22	0/1/1/1
3	NAG	F	1	1,3	-	1/6/23/26	0/1/1/1
3	NAG	F	2	3	-	0/6/23/26	0/1/1/1
3	BMA	F	3	3	-	2/2/19/22	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	-	0/2/19/22	0/1/1/1
4	NAG	G	1	1,4	-	3/6/23/26	0/1/1/1
4	NAG	G	2	4	-	5/6/23/26	0/1/1/1
5	Z3Q	H	1	5	-	9/12/32/32	0/1/1/1
5	GAL	H	2	5	-	1/2/19/22	0/1/1/1
5	5N6	H	3	5	-	10/17/37/41	0/1/1/1
4	NAG	I	1	2,4	-	0/6/23/26	0/1/1/1
4	NAG	I	2	4	-	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

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
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	-	2/2/19/22	0/1/1/1
3	MAN	J	6	3	-	0/2/19/22	0/1/1/1
4	NAG	K	1	1,4	-	0/6/23/26	0/1/1/1
4	NAG	K	2	4	-	3/6/23/26	0/1/1/1
4	NAG	L	1	1,4	-	1/6/23/26	0/1/1/1
4	NAG	L	2	4	-	2/6/23/26	0/1/1/1
5	Z3Q	M	1	5	-	5/12/32/32	0/1/1/1
5	GAL	M	2	5	-	2/2/19/22	0/1/1/1
5	5N6	M	3	5	-	1/17/37/41	0/1/1/1
4	NAG	N	1	2,4	-	0/6/23/26	0/1/1/1
4	NAG	N	2	4	-	0/6/23/26	0/1/1/1

All (21) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	H	1	Z3Q	N1B-N1A	-5.98	1.07	1.23
5	M	3	5N6	C4-C5	-5.40	1.48	1.53
5	M	3	5N6	O4-C4	-4.78	1.33	1.43
5	H	3	5N6	C4-C5	-4.67	1.49	1.53
5	M	3	5N6	C7-C6	-4.15	1.47	1.53
5	M	1	Z3Q	N1B-N1A	-4.11	1.12	1.23
5	M	3	5N6	O9-C9	-3.30	1.37	1.45
5	H	1	Z3Q	O1-C1	3.30	1.45	1.40
5	H	3	5N6	O4-C4	-3.28	1.36	1.43
5	M	3	5N6	C8-C7	-3.00	1.47	1.53
5	H	3	5N6	O9-C9	-2.70	1.39	1.45
5	M	3	5N6	O6-C6	-2.66	1.39	1.44
3	E	3	BMA	O5-C1	-2.66	1.39	1.43
5	M	3	5N6	C3-C4	-2.58	1.48	1.52
5	H	3	5N6	C9-C8	-2.53	1.48	1.51
5	M	3	5N6	O8-C8	-2.46	1.38	1.43
5	H	3	5N6	C3-C4	-2.37	1.48	1.52
5	M	3	5N6	O6-C2	-2.22	1.38	1.43
5	M	1	Z3Q	O1-C1	2.22	1.44	1.40
5	H	2	GAL	O5-C5	-2.07	1.39	1.43
4	G	1	NAG	C1-C2	2.04	1.55	1.52

All (74) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	H	1	Z3Q	C1-C2-N2	-12.20	90.00	111.00
5	H	1	Z3Q	O5-C1-O1	-11.54	82.64	109.97
3	F	3	BMA	C2-C3-C4	-8.31	96.51	110.89
3	E	3	BMA	C1-O5-C5	-7.23	102.39	112.19
3	E	3	BMA	O2-C2-C3	-6.14	97.83	110.14
3	F	3	BMA	C1-C2-C3	6.04	117.09	109.67
3	F	4	MAN	C1-O5-C5	5.47	119.60	112.19
3	F	2	NAG	O4-C4-C3	4.64	121.07	110.35
3	F	4	MAN	C1-C2-C3	4.24	114.88	109.67
5	H	2	GAL	C1-O5-C5	-4.20	106.50	112.19
5	H	2	GAL	O5-C5-C6	4.08	113.60	107.20
4	G	2	NAG	C1-O5-C5	3.85	117.41	112.19
5	H	1	Z3Q	C3-C2-N2	-3.81	103.42	110.62
3	J	4	MAN	C1-O5-C5	3.80	117.34	112.19
3	E	5	MAN	O5-C5-C6	3.76	113.10	107.20
5	H	3	5N6	C9-O9-CAG	-3.75	107.69	117.10
5	H	1	Z3Q	C11-O1-C1	3.73	120.03	113.84
5	H	2	GAL	O3-C3-C4	-3.70	101.80	110.35
4	I	2	NAG	C1-O5-C5	3.67	117.17	112.19
4	G	1	NAG	C1-O5-C5	3.61	117.08	112.19
5	H	1	Z3Q	C2-N2-C7	3.58	131.88	123.18
4	N	2	NAG	C1-O5-C5	3.51	116.95	112.19
3	F	3	BMA	O4-C4-C3	3.40	118.22	110.35
5	H	3	5N6	C9-C8-C7	-3.39	105.65	112.20
4	L	2	NAG	C1-O5-C5	3.28	116.63	112.19
5	H	1	Z3Q	O5-C5-C6	-3.23	98.41	106.44
5	H	2	GAL	C2-C3-C4	3.04	116.16	110.89
4	N	2	NAG	C2-N2-C7	-3.00	118.64	122.90
3	F	3	BMA	O6-C6-C5	-2.98	101.05	111.29
4	L	1	NAG	C1-O5-C5	2.92	116.15	112.19
3	F	3	BMA	O3-C3-C4	2.90	117.06	110.35
3	E	5	MAN	C6-C5-C4	-2.89	106.24	113.00
4	I	2	NAG	C2-N2-C7	-2.75	118.99	122.90
3	F	3	BMA	O3-C3-C2	-2.75	104.73	109.99
3	F	6	MAN	O5-C5-C6	2.74	111.50	107.20
3	F	2	NAG	O4-C4-C5	-2.73	102.52	109.30
3	E	6	MAN	O5-C5-C6	2.69	111.42	107.20
3	F	2	NAG	O3-C3-C2	-2.67	103.95	109.47
5	M	1	Z3Q	C11-O1-C1	-2.63	109.48	113.84
5	M	1	Z3Q	C1-O5-C5	-2.62	108.55	113.69
3	J	2	NAG	O4-C4-C5	-2.61	102.83	109.30
5	H	3	5N6	C4-C5-N5	-2.58	105.27	110.38
3	E	3	BMA	O6-C6-C5	-2.47	102.81	111.29

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	H	3	5N6	O10-C10-N5	2.45	126.46	121.95
5	M	2	GAL	C2-C3-C4	2.45	115.13	110.89
5	H	1	Z3Q	O4-C4-C5	-2.44	103.24	109.30
3	F	1	NAG	O5-C5-C6	2.43	111.02	107.20
3	F	2	NAG	C1-O5-C5	2.43	115.48	112.19
5	M	2	GAL	O5-C5-C4	-2.40	104.99	110.83
4	K	1	NAG	C2-N2-C7	-2.39	119.50	122.90
4	N	1	NAG	C1-O5-C5	2.37	115.40	112.19
5	M	3	5N6	C9-C8-C7	-2.33	107.70	112.20
3	J	4	MAN	C6-C5-C4	-2.32	107.56	113.00
3	J	5	MAN	O5-C5-C6	2.30	110.82	107.20
3	E	4	MAN	C1-C2-C3	2.30	112.50	109.67
3	F	4	MAN	O3-C3-C4	-2.30	105.04	110.35
5	H	3	5N6	C6-O6-C2	2.26	116.17	111.34
4	K	1	NAG	C1-O5-C5	2.25	115.24	112.19
3	J	4	MAN	C1-C2-C3	2.24	112.42	109.67
3	J	4	MAN	O3-C3-C4	-2.22	105.21	110.35
4	G	2	NAG	C1-C2-N2	-2.22	106.70	110.49
4	G	1	NAG	O5-C5-C6	2.21	110.67	107.20
5	H	3	5N6	O8-C8-C9	-2.17	105.03	109.92
5	H	3	5N6	C4-C5-C6	-2.16	103.63	109.10
3	E	3	BMA	O5-C1-C2	-2.15	107.45	110.77
3	E	3	BMA	O5-C5-C4	-2.15	105.61	110.83
5	M	2	GAL	O3-C3-C4	-2.12	105.44	110.35
4	G	2	NAG	C2-N2-C7	2.10	125.89	122.90
3	J	3	BMA	O3-C3-C2	2.10	114.01	109.99
5	H	3	5N6	O9-C9-C8	-2.06	103.81	108.45
4	N	1	NAG	O4-C4-C5	-2.05	104.20	109.30
3	F	5	MAN	O5-C5-C6	2.02	110.37	107.20
3	E	1	NAG	C1-O5-C5	2.01	114.92	112.19
3	E	1	NAG	C2-N2-C7	-2.01	120.04	122.90

There are no chirality outliers.

All (49) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	H	3	5N6	C7-C8-C9-O9
5	H	3	5N6	O8-C8-C9-O9
5	H	3	5N6	O6-C6-C7-C8
5	H	3	5N6	O6-C6-C7-O7
5	H	1	Z3Q	C2-C1-O1-C11
5	H	1	Z3Q	O5-C1-O1-C11

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Mol	Chain	Res	Type	Atoms
5	H	1	Z3Q	O1-C11-C12-N1A
5	H	1	Z3Q	C1-C2-N2-C7
4	K	2	NAG	C8-C7-N2-C2
4	K	2	NAG	O7-C7-N2-C2
5	M	1	Z3Q	C2-C1-O1-C11
5	H	3	5N6	CAF-CAG-O9-C9
4	G	1	NAG	O5-C5-C6-O6
3	F	3	BMA	O5-C5-C6-O6
5	M	1	Z3Q	O5-C5-C6-O6
5	M	2	GAL	O5-C5-C6-O6
4	L	2	NAG	O5-C5-C6-O6
3	F	3	BMA	C4-C5-C6-O6
5	H	1	Z3Q	O5-C5-C6-O6
4	G	2	NAG	O5-C5-C6-O6
5	M	2	GAL	C4-C5-C6-O6
4	G	1	NAG	C4-C5-C6-O6
5	M	1	Z3Q	C4-C5-C6-O6
5	H	3	5N6	OBJ-CAG-O9-C9
5	H	1	Z3Q	C8-C7-N2-C2
5	H	1	Z3Q	O7-C7-N2-C2
4	G	2	NAG	C8-C7-N2-C2
4	G	2	NAG	C4-C5-C6-O6
4	L	2	NAG	C4-C5-C6-O6
4	G	2	NAG	O7-C7-N2-C2
5	H	1	Z3Q	C4-C5-C6-O6
3	J	5	MAN	O5-C5-C6-O6
3	J	5	MAN	C4-C5-C6-O6
3	E	5	MAN	O5-C5-C6-O6
5	H	3	5N6	C4-C5-N5-C10
5	M	1	Z3Q	O5-C1-O1-C11
3	E	5	MAN	C4-C5-C6-O6
5	M	1	Z3Q	C12-N1A-N1B-N1C
3	F	1	NAG	C4-C5-C6-O6
5	H	3	5N6	C5-C6-C7-C8
5	H	3	5N6	C5-C6-C7-O7
5	H	2	GAL	C4-C5-C6-O6
5	H	1	Z3Q	C3-C2-N2-C7
4	K	2	NAG	C4-C5-C6-O6
5	H	3	5N6	C6-C5-N5-C10
5	M	3	5N6	CAF-CAG-O9-C9
4	G	2	NAG	C1-C2-N2-C7
4	G	1	NAG	C8-C7-N2-C2

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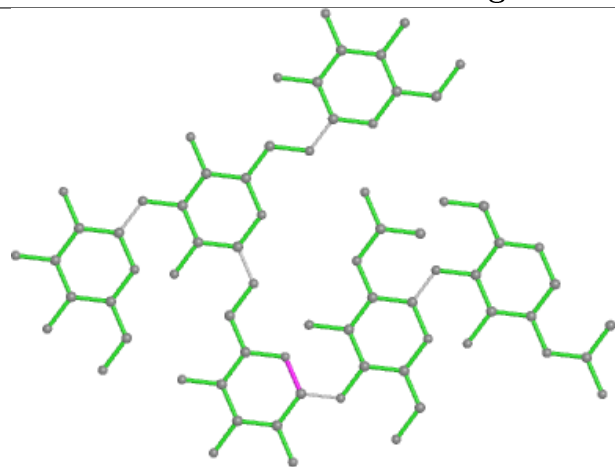
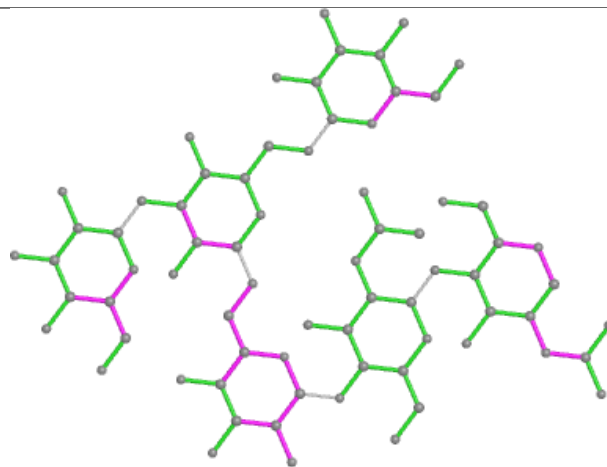
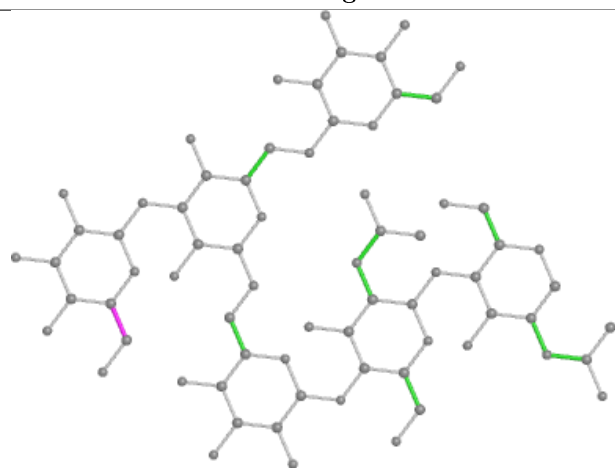
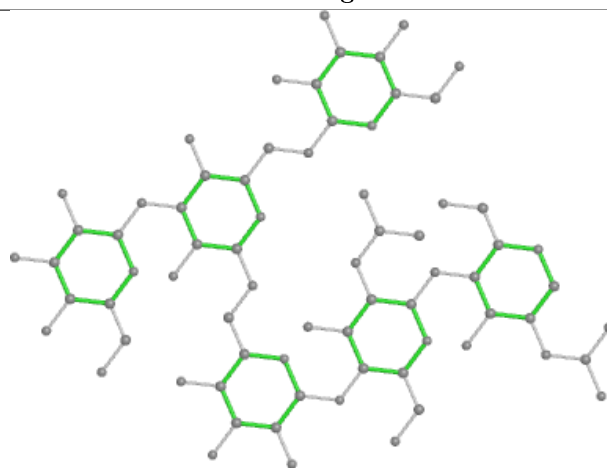
Mol	Chain	Res	Type	Atoms
4	L	1	NAG	C4-C5-C6-O6

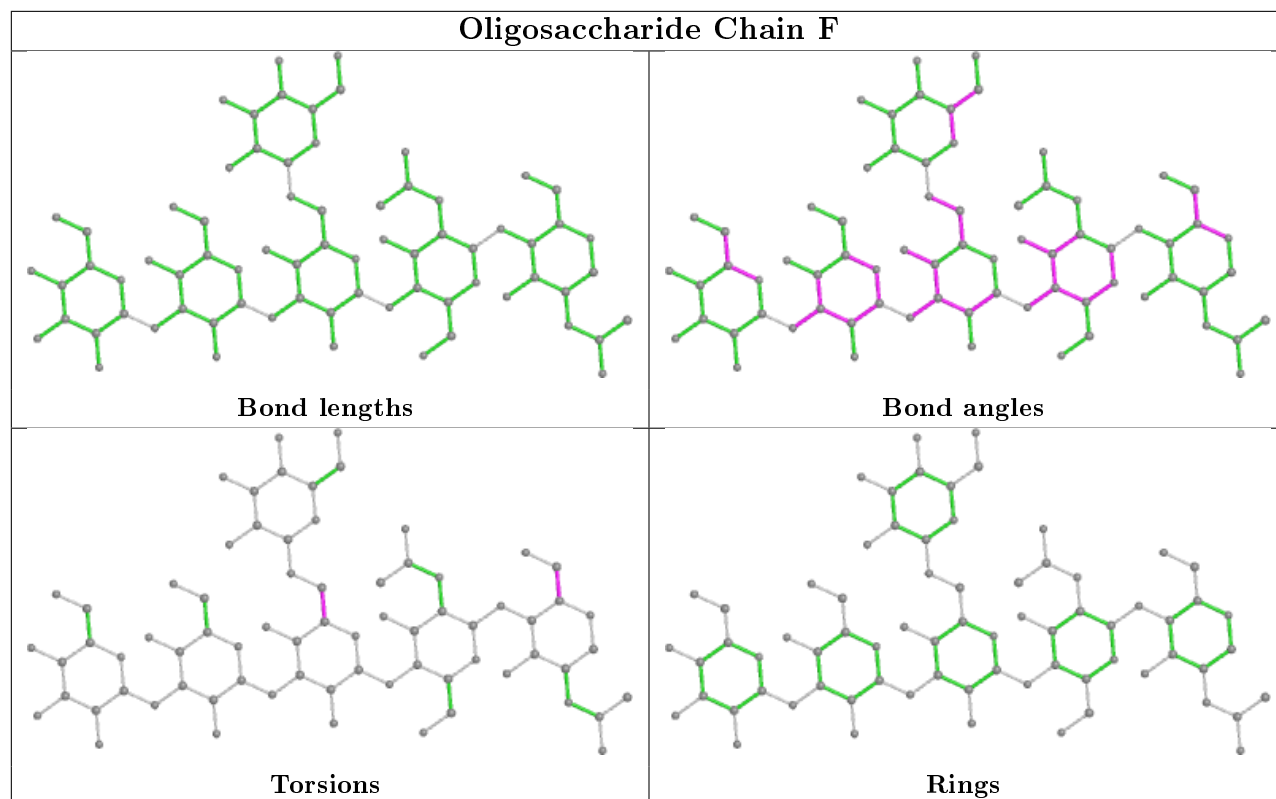
There are no ring outliers.

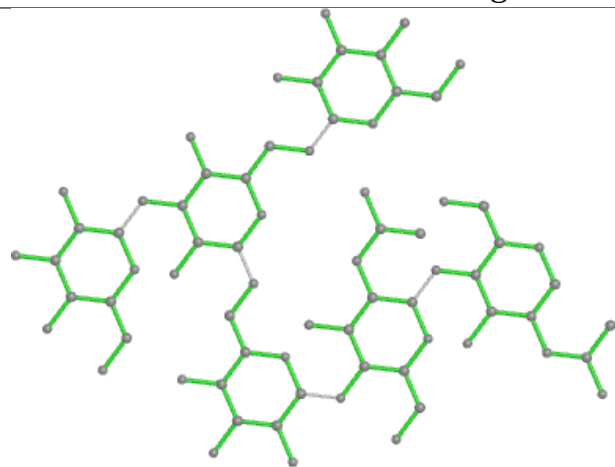
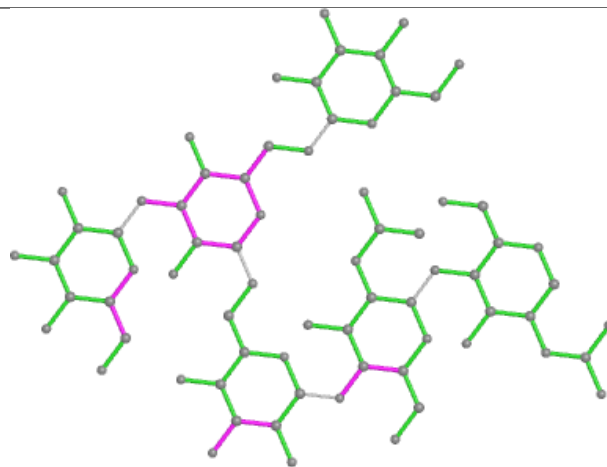
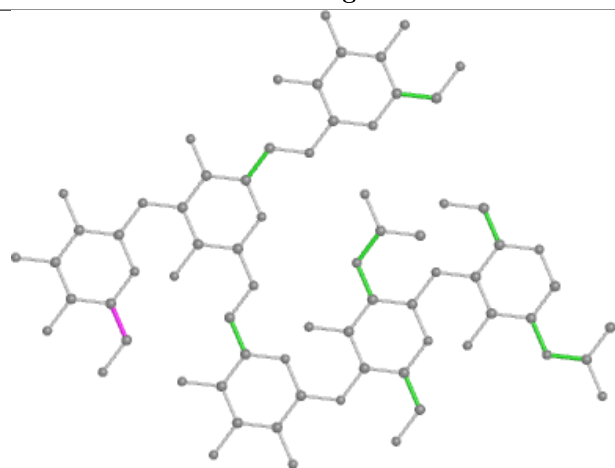
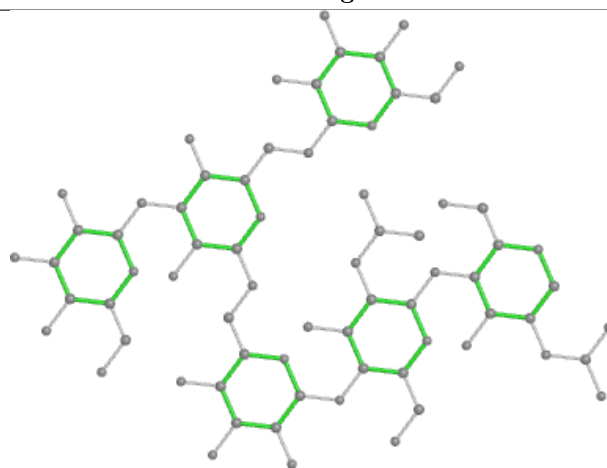
13 monomers are involved in 15 short contacts:

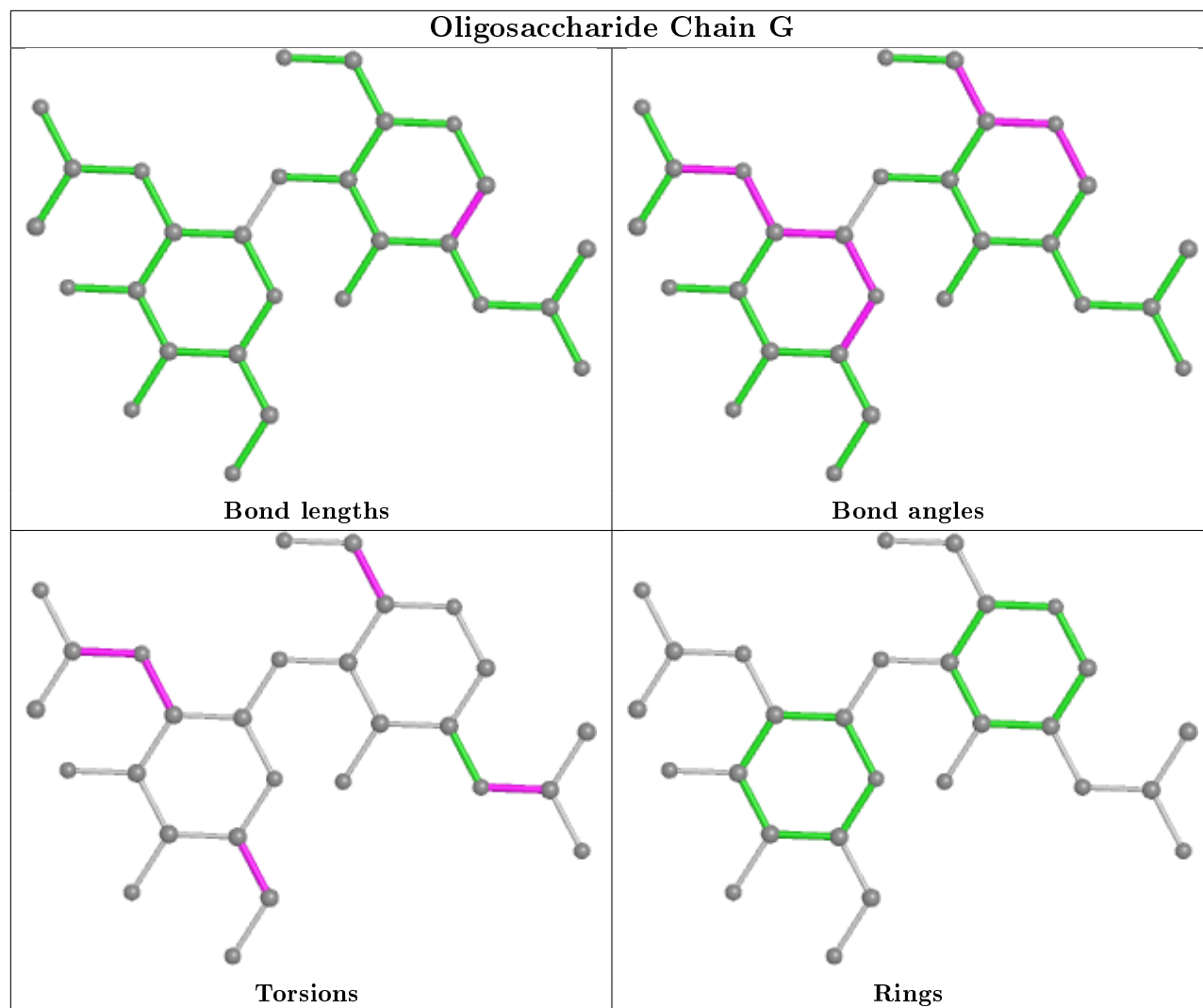
Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	M	3	5N6	2	0
5	H	2	GAL	3	0
3	F	2	NAG	1	0
3	F	3	BMA	1	0
3	E	3	BMA	1	0
5	H	3	5N6	5	0
3	E	2	NAG	1	0
4	N	2	NAG	1	0
5	H	1	Z3Q	1	0
4	I	2	NAG	2	0
4	L	1	NAG	1	0
5	M	2	GAL	2	0
4	G	1	NAG	1	0

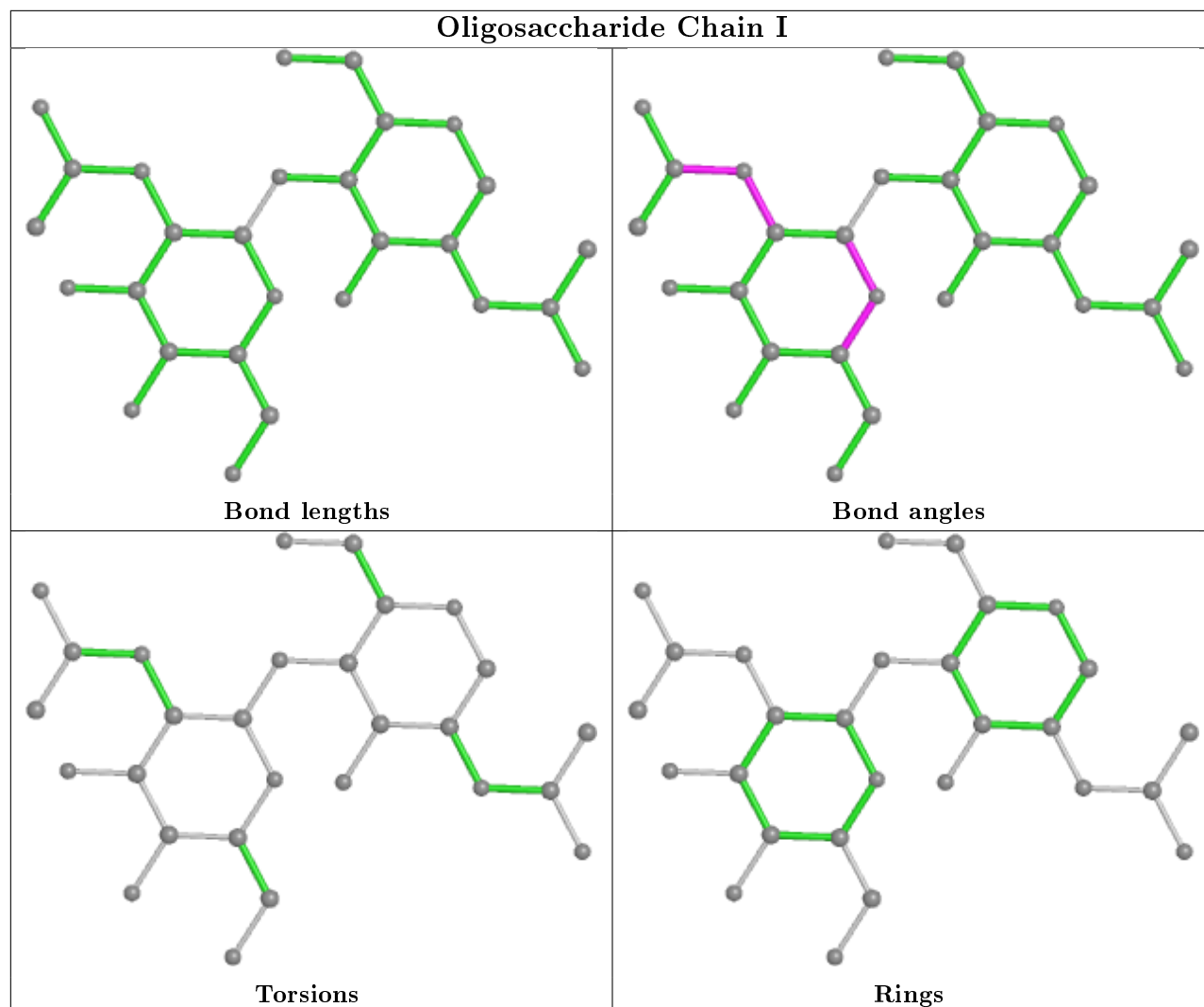
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for oligosaccharide.

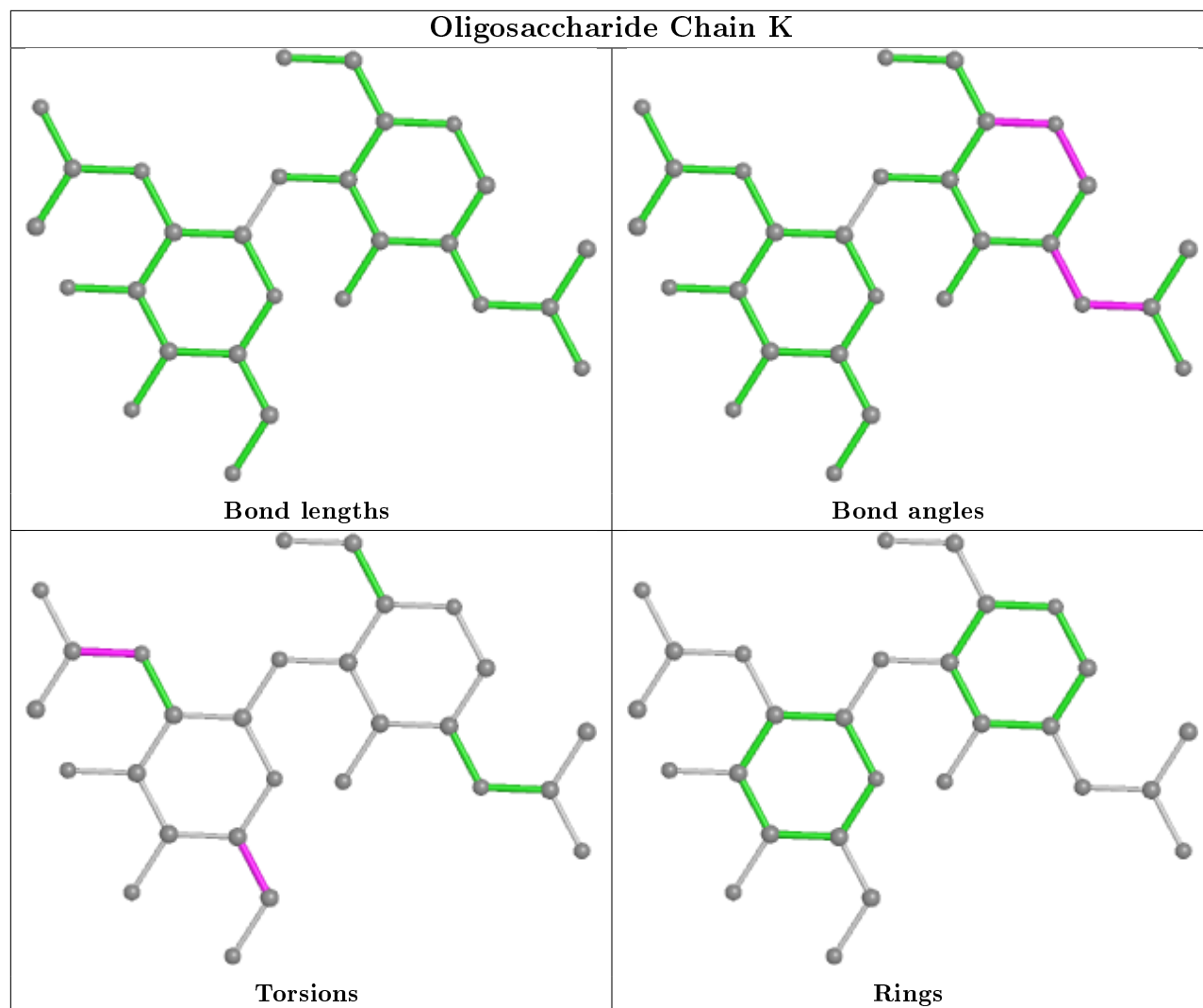
Oligosaccharide Chain E**Bond lengths****Bond angles****Torsions****Rings**

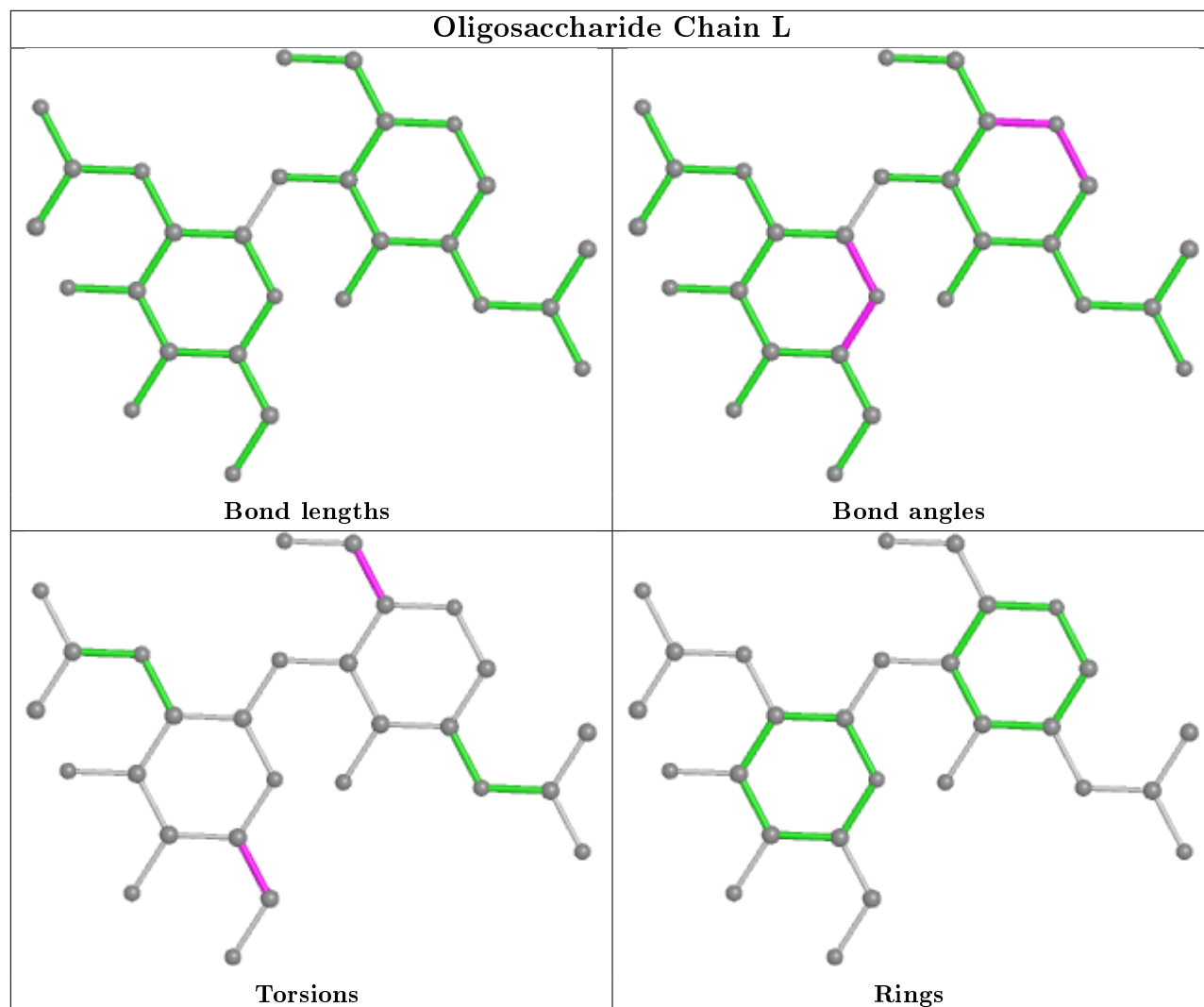


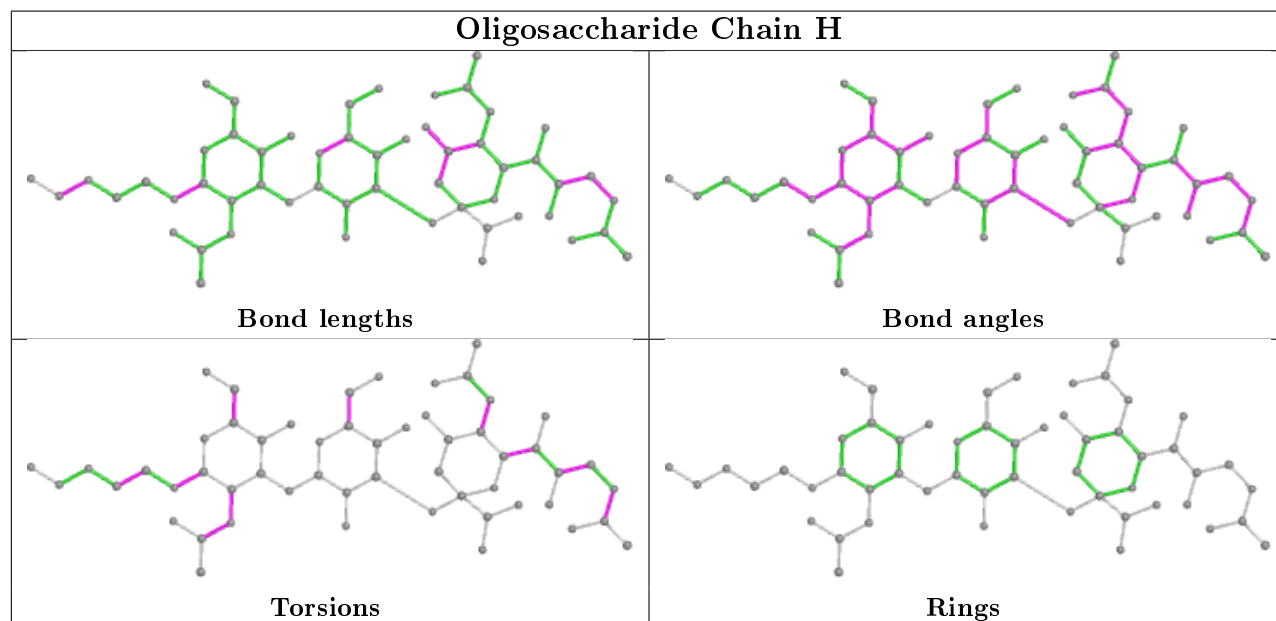
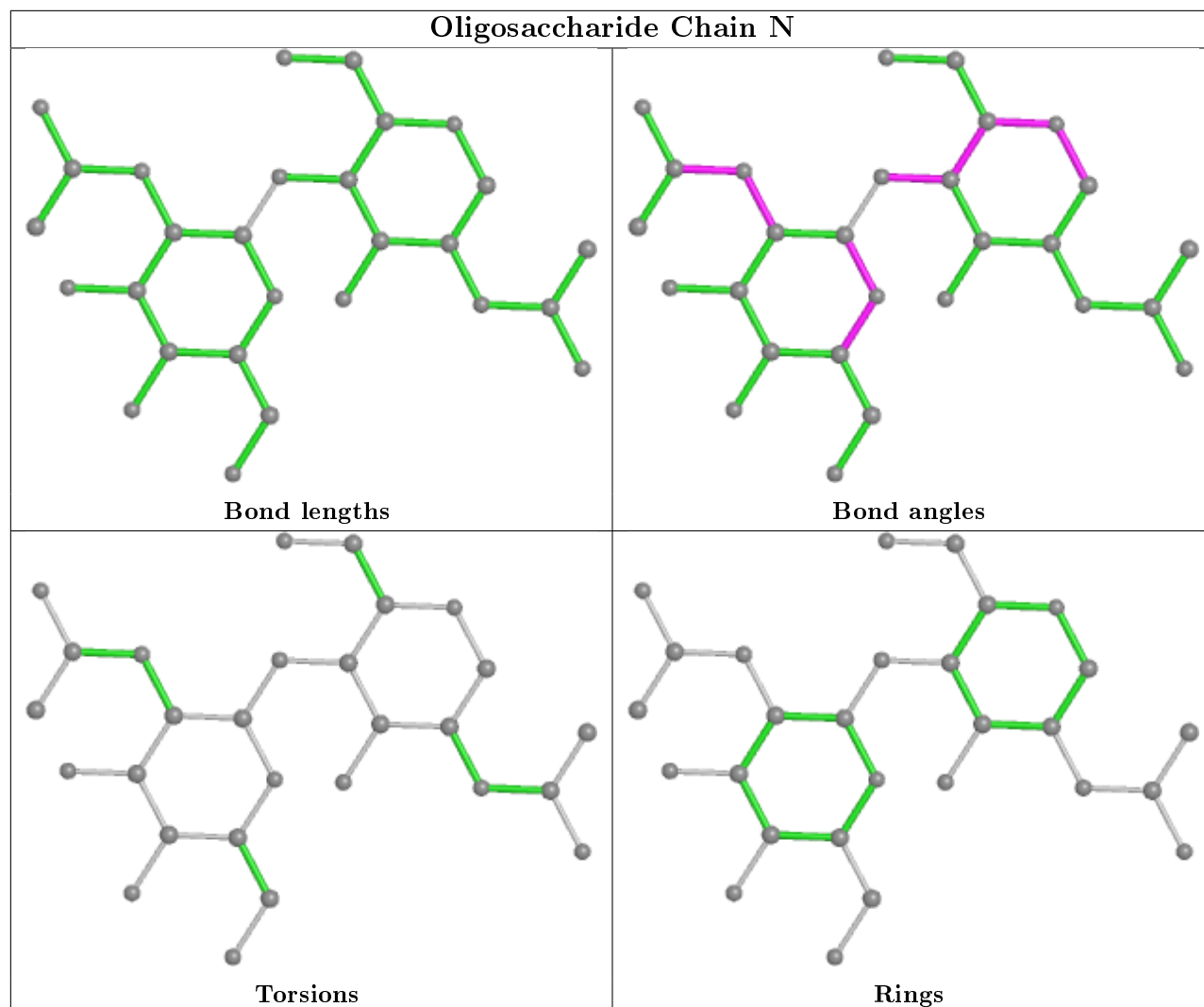
Oligosaccharide Chain J**Bond lengths****Bond angles****Torsions****Rings**

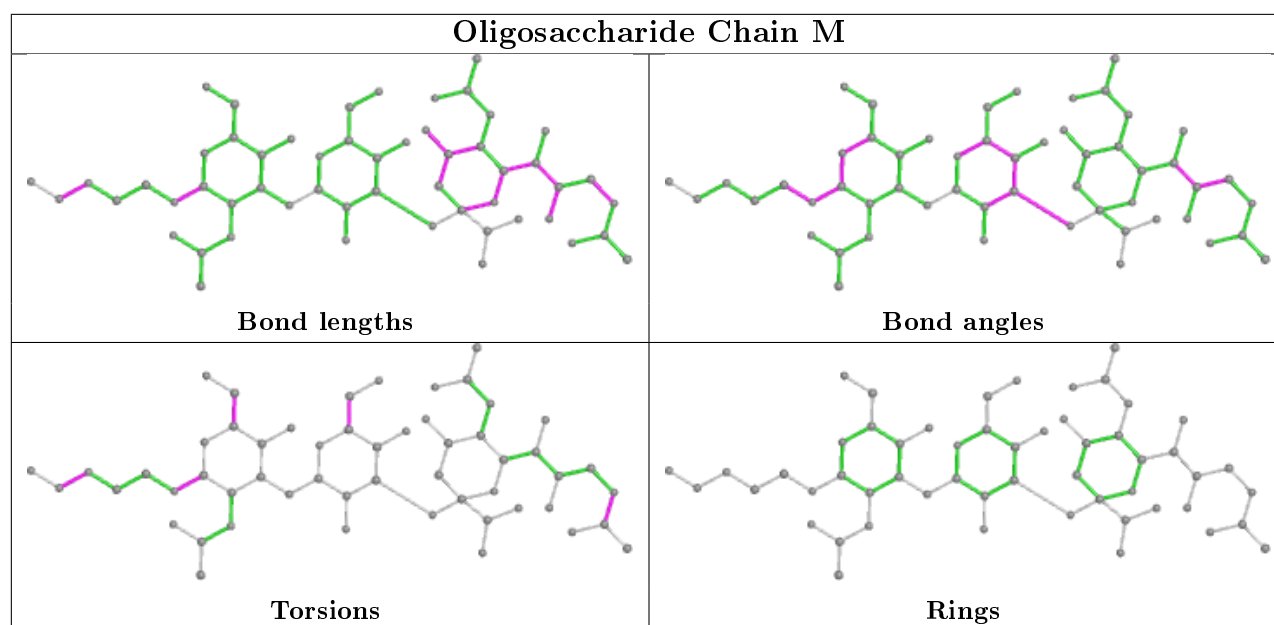












5.6 Ligand geometry [i](#)

2 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
6	NAG	D	703	2	14,14,15	0.32	0	17,19,21	1.38	2 (11%)
6	NAG	B	703	2	14,14,15	0.38	0	17,19,21	1.38	3 (17%)

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	D	703	2	-	4/6/23/26	0/1/1/1
6	NAG	B	703	2	-	1/6/23/26	0/1/1/1

There are no bond length outliers.

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	D	703	NAG	C1-O5-C5	4.00	117.62	112.19
6	B	703	NAG	C1-O5-C5	3.57	117.03	112.19
6	B	703	NAG	C4-C3-C2	-2.51	107.34	111.02
6	D	703	NAG	C2-N2-C7	-2.30	119.63	122.90
6	B	703	NAG	C2-N2-C7	-2.04	120.00	122.90

There are no chirality outliers.

All (5) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
6	D	703	NAG	C4-C5-C6-O6
6	D	703	NAG	C8-C7-N2-C2
6	D	703	NAG	O5-C5-C6-O6
6	D	703	NAG	O7-C7-N2-C2
6	B	703	NAG	C4-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 ⓘ

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, 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	424/427 (99%)	-0.18	20 (4%) 31 30	28, 40, 98, 144	0
1	C	424/427 (99%)	-0.48	2 (0%) 91 90	24, 36, 60, 78	0
2	B	149/149 (100%)	0.52	20 (13%) 3 2	26, 55, 120, 198	0
2	D	149/149 (100%)	0.53	13 (8%) 10 8	30, 51, 127, 190	0
All	All	1146/1152 (99%)	-0.11	55 (4%) 30 29	24, 40, 98, 198	0

All (55) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	D	32	GLY	17.8
2	B	34	GLY	14.7
2	B	37	ALA	14.1
2	D	33	GLY	13.8
2	D	37	ALA	11.0
2	B	32	GLY	10.5
2	D	35	GLY	10.0
1	A	191	VAL	8.5
1	A	193	PRO	7.8
2	D	36	GLY	7.1
2	B	33	GLY	6.9
2	D	34	GLY	6.2
1	A	237	PHE	6.0
1	A	264	VAL	5.5
2	B	36	GLY	5.4
2	B	44	ALA	5.3
2	D	149	SER	4.6
1	A	265	ALA	4.4
2	D	143	SER	4.3
2	D	31	ASN	4.2
1	A	199	VAL	4.0

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Mol	Chain	Res	Type	RSRZ
2	B	141	VAL	3.8
2	B	45	GLY	3.7
2	D	38	SER	3.7
2	D	30	SER	3.6
1	A	241	GLY	3.5
2	B	157	PHE	3.5
2	B	30	SER	3.2
1	A	273	ALA	3.2
1	A	236	LYS	3.1
2	B	131	ALA	3.1
1	A	197	PRO	3.1
1	A	267	TYR	3.0
1	C	270	ILE	2.9
1	A	238	ASN	2.8
1	A	235	ALA	2.8
1	A	240	PHE	2.7
2	B	38	SER	2.7
1	A	192	ASN	2.6
2	B	147	ASP	2.5
2	B	149	SER	2.5
2	D	142	ASP	2.5
2	B	31	ASN	2.5
2	B	9	PHE	2.5
1	A	296	ARG	2.4
1	A	127	PHE	2.3
1	C	236	LYS	2.2
2	B	145	ASN	2.2
2	D	45	GLY	2.2
1	A	198	GLN	2.2
2	B	29	ARG	2.2
2	B	16	PHE	2.1
2	B	49	ILE	2.0
1	A	275	VAL	2.0
1	A	10	ARG	2.0

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 ⓘ

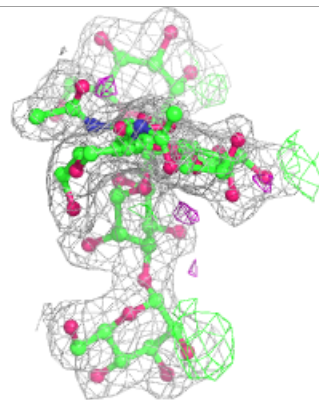
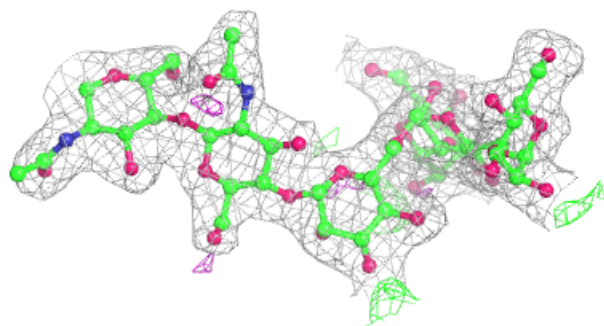
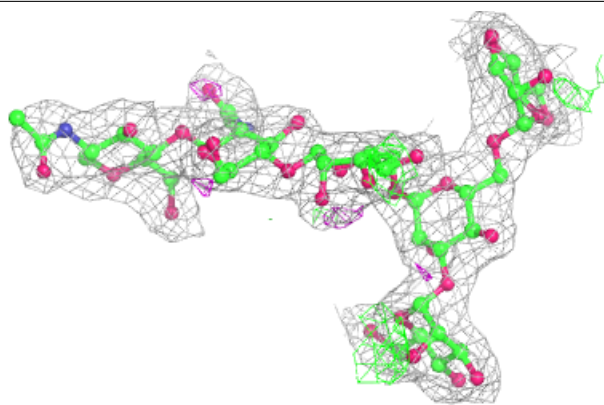
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q<0.9' lists 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	Z3Q	M	1	20/20	0.45	0.66	109,126,133,138	0
5	Z3Q	H	1	20/20	0.61	0.62	129,139,148,148	0
5	GAL	H	2	11/12	0.77	0.31	96,109,120,121	0
4	NAG	G	2	14/15	0.81	0.13	58,65,73,79	0
5	GAL	M	2	11/12	0.86	0.23	68,84,101,102	0
3	BMA	F	3	11/12	0.88	0.17	43,47,54,60	0
3	MAN	F	5	11/12	0.89	0.16	52,58,65,65	0
4	NAG	L	2	14/15	0.90	0.10	53,57,63,68	0
4	NAG	K	2	14/15	0.91	0.14	39,46,51,52	0
3	MAN	E	5	11/12	0.92	0.09	32,38,42,46	0
3	BMA	J	3	11/12	0.93	0.09	33,35,39,41	0
3	MAN	E	6	11/12	0.93	0.09	37,40,43,44	0
3	NAG	E	2	14/15	0.93	0.12	36,42,46,50	0
3	MAN	F	6	11/12	0.93	0.10	39,42,47,47	0
3	BMA	E	3	11/12	0.94	0.10	35,38,42,43	0
4	NAG	I	2	14/15	0.94	0.12	43,48,56,58	0
4	NAG	N	2	14/15	0.94	0.12	36,44,50,54	0
3	MAN	F	4	11/12	0.94	0.11	42,47,48,51	0
3	MAN	J	5	11/12	0.94	0.09	38,40,45,48	0
5	5N6	H	3	23/24	0.94	0.19	71,81,85,87	0
4	NAG	I	1	14/15	0.94	0.10	30,35,39,41	0
4	NAG	G	1	14/15	0.94	0.09	60,63,69,72	0
3	MAN	J	6	11/12	0.94	0.09	41,46,52,59	0
3	NAG	E	1	14/15	0.95	0.09	38,40,49,50	0
3	NAG	F	2	14/15	0.95	0.10	31,36,38,39	0
5	5N6	M	3	23/24	0.96	0.13	43,52,60,62	0
3	NAG	J	1	14/15	0.96	0.08	34,37,41,42	0
3	NAG	J	2	14/15	0.96	0.09	28,33,38,40	0
4	NAG	L	1	14/15	0.96	0.10	51,52,64,68	0
3	NAG	F	1	14/15	0.96	0.08	29,37,40,42	0
4	NAG	N	1	14/15	0.96	0.09	29,34,36,36	0
3	MAN	J	4	11/12	0.96	0.07	30,34,39,41	0
3	MAN	E	4	11/12	0.97	0.09	28,32,37,38	0
4	NAG	K	1	14/15	0.98	0.08	25,30,36,39	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.

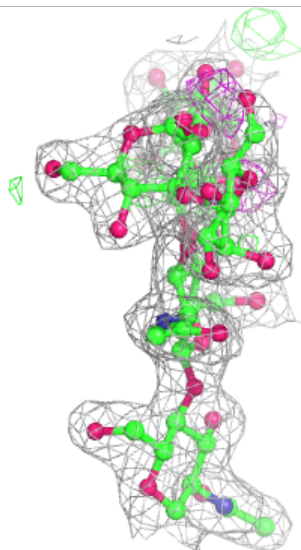
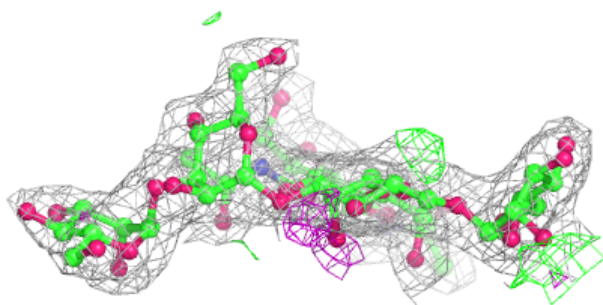
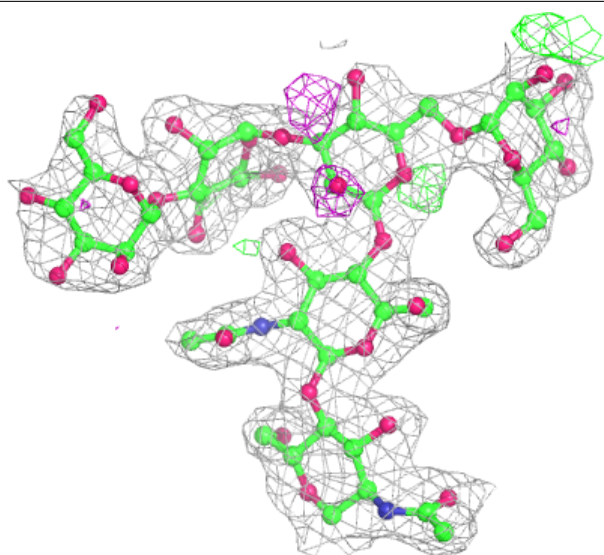
Electron density around Chain E:

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



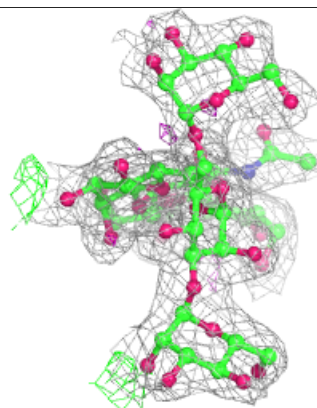
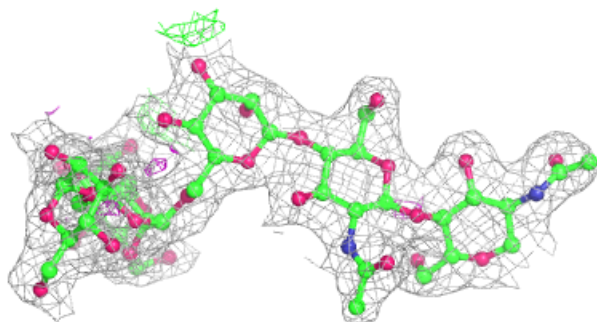
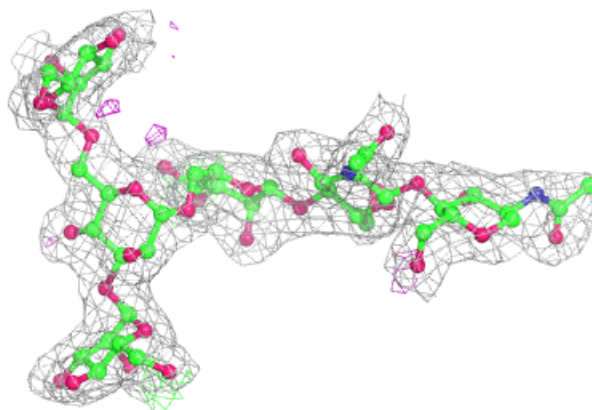
Electron density around Chain F:

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 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



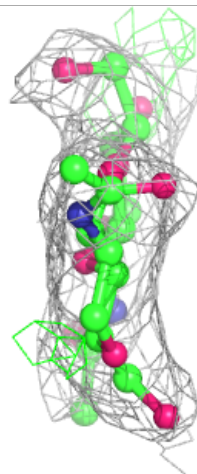
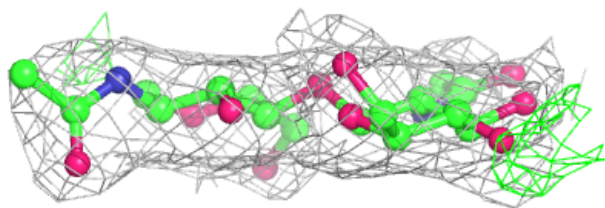
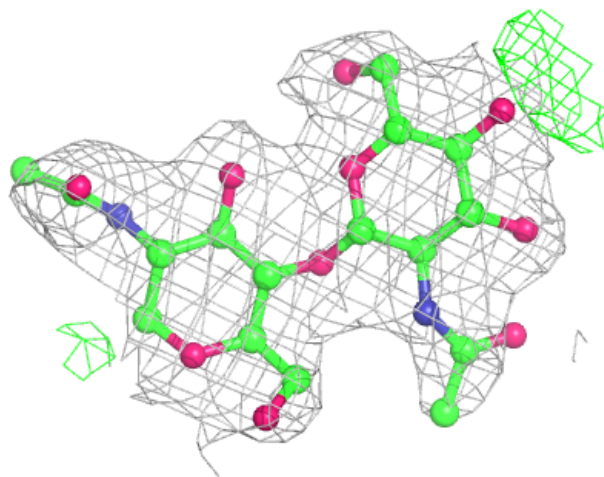
Electron density around Chain J:

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 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



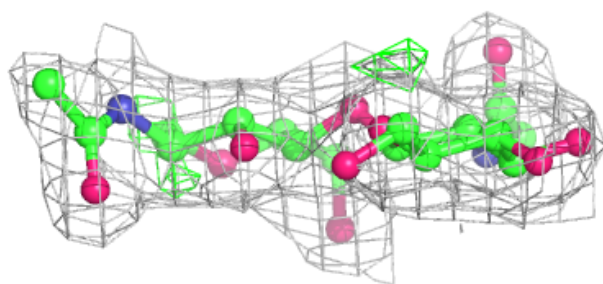
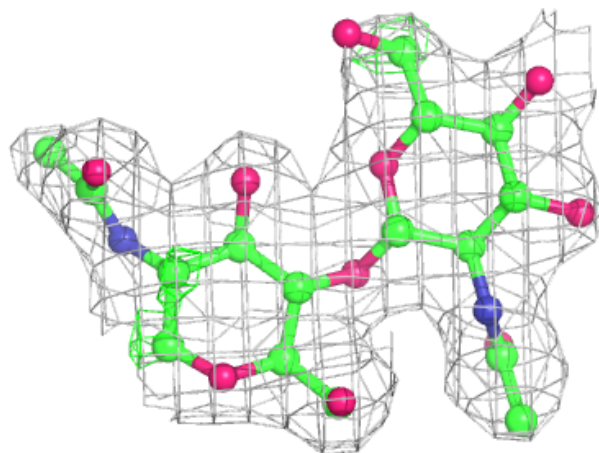
Electron density around Chain G:

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



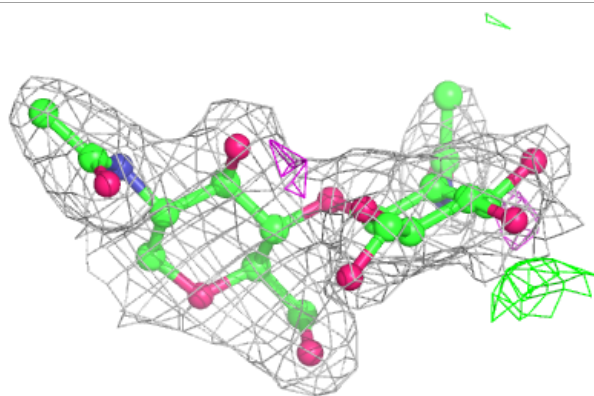
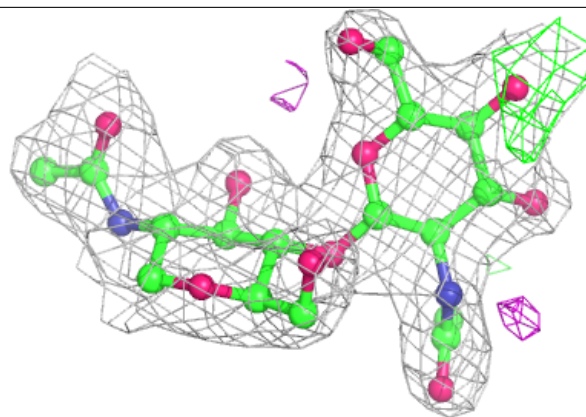
Electron density around Chain I:

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

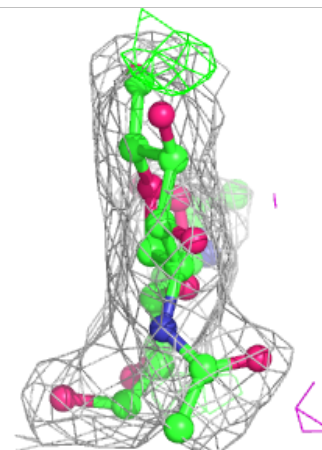
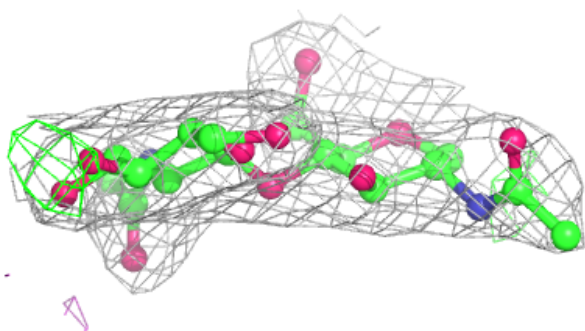
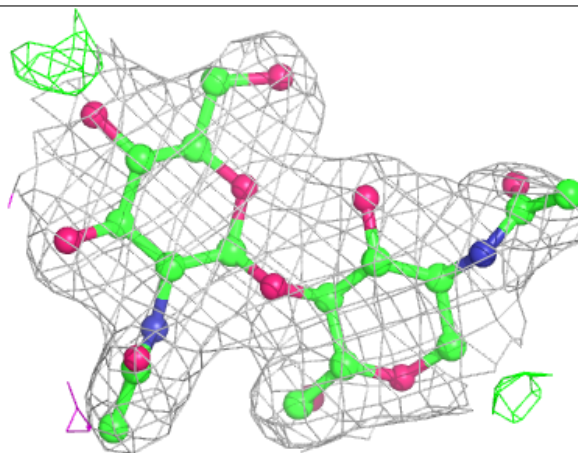


Electron density around Chain K:

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

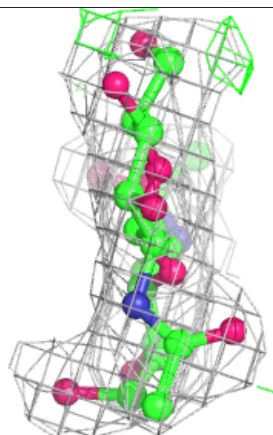
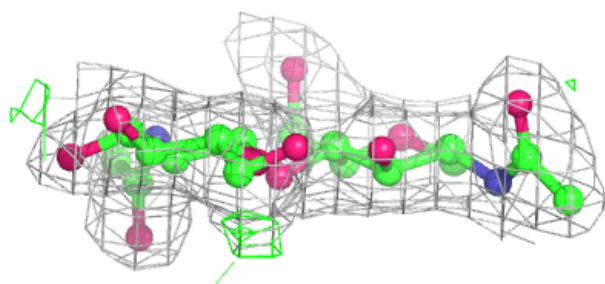
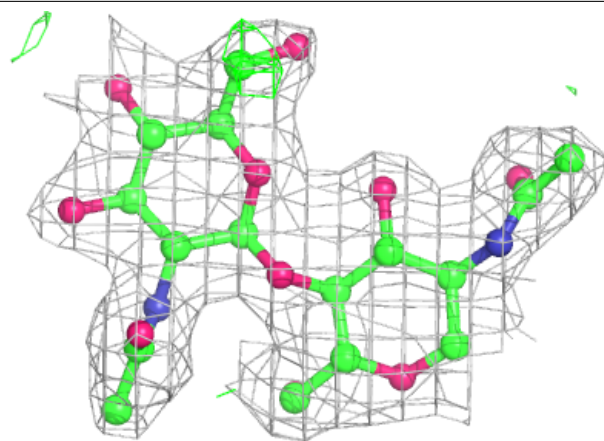
**Electron density around Chain L:**

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

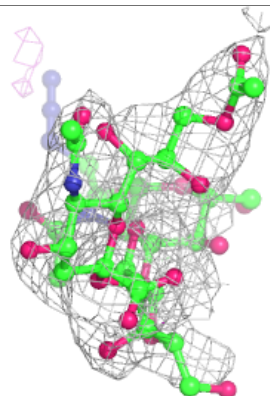
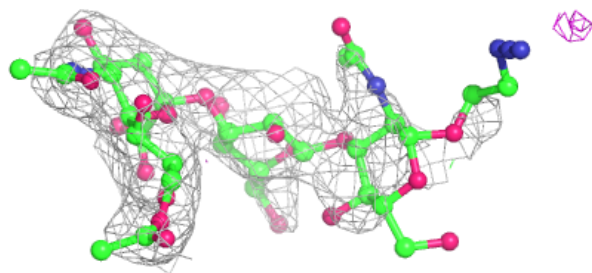
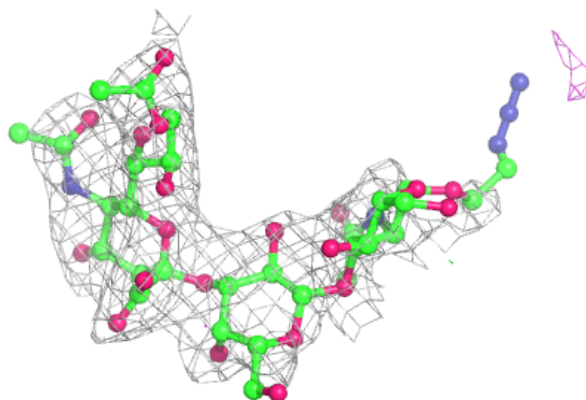


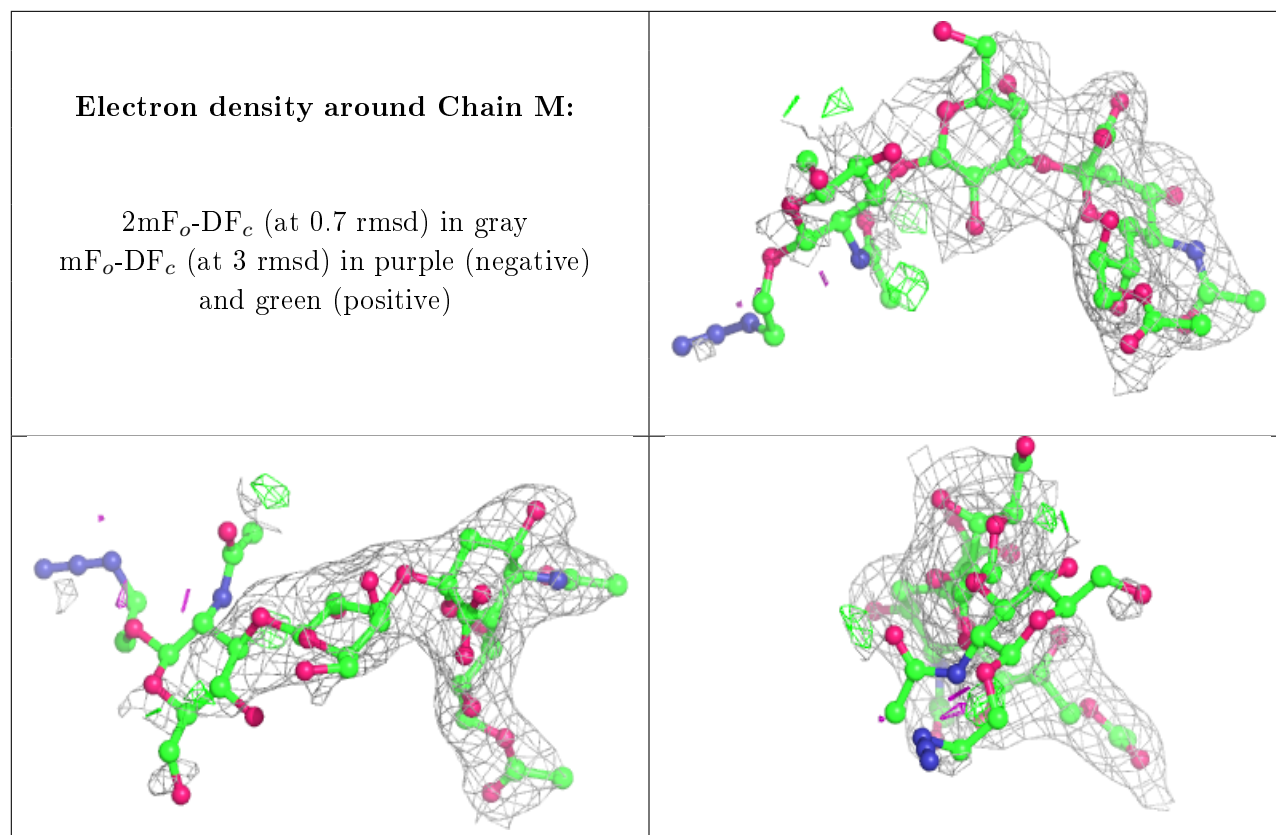
Electron density around Chain N:

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

**Electron density around Chain H:**

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





6.4 Ligands [i](#)

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
6	NAG	D	703	14/15	0.79	0.25	88,99,103,103	0
6	NAG	B	703	14/15	0.86	0.21	70,77,81,82	0

6.5 Other polymers [i](#)

There are no such residues in this entry.