Full wwPDB X-ray Structure Validation Report (i)

Jun 15, 2024 – 09:35 AM EDT

PDB ID	:	2Q9O
Title	:	Near-atomic resolution structure of a Melanocarpus albomyces laccase
Authors	:	Hakulinen, N.; Rouvinen, J.
Deposited on	:	2007-06-13
Resolution	:	1.30 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 (i) symbol.

The types of validation reports are described at
http://www.wwpdb.org/validation/2017/FAQs#types.

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

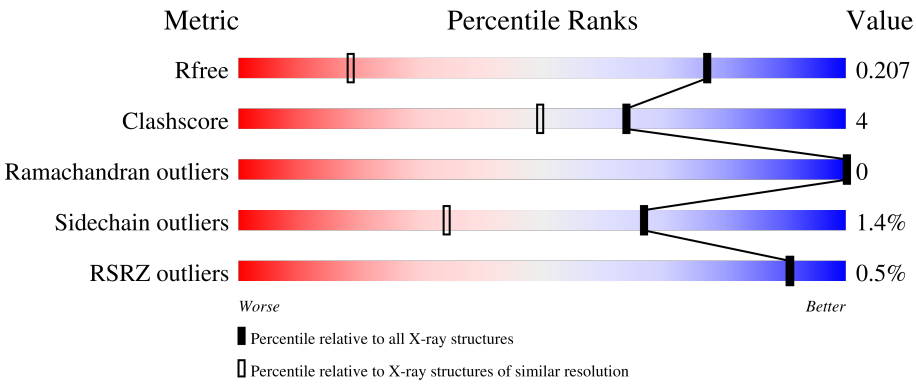
MolProbity	:	4.02b-467
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	1.20.1
EDS	:	2.37.1
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.37.1

1 Overall quality at a glance (i)

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

The reported resolution of this entry is 1.30 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	1058 (1.30-1.30)
Clashscore	141614	1101 (1.30-1.30)
Ramachandran outliers	138981	1058 (1.30-1.30)
Sidechain outliers	138945	1058 (1.30-1.30)
RSRZ outliers	127900	1029 (1.30-1.30)

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

Mol	Chain	Length	Quality of chain
1	A	559	% 88% 11% •
1	B	559	90% 9% •
2	C	2	100%
2	E	2	100%
2	I	2	100%

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Mol	Chain	Length	Quality of chain
2	J	2	100%
2	K	2	100%
3	D	3	100%
3	F	3	100%
3	H	3	100%
4	G	5	20% 60% 20%
4	L	5	20% 80%

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
3	MAN	D	3	X	-	-	-
3	MAN	F	3	X	-	-	-
3	MAN	H	3	X	-	-	-
4	MAN	G	3	X	-	-	-
4	MAN	L	3	X	-	-	-

2 Entry composition (i)

There are 11 unique types of molecules in this entry. The entry contains 10805 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 Laccase-1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	559	4432	2798	770	847	17	0	14	0
1	B	559	4418	2789	767	846	16	0	15	0

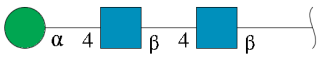
There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	98	OHI	HIS	modified residue	UNP Q70KY3
B	98	OHI	HIS	modified residue	UNP Q70KY3

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

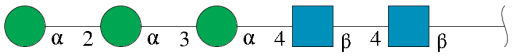
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
			Total	C	N	O			
2	C	2	28	16	2	10	0	0	0
2	E	2	28	16	2	10	0	0	0
2	I	2	28	16	2	10	0	0	0
2	J	2	28	16	2	10	0	0	0
2	K	2	28	16	2	10	0	0	0

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



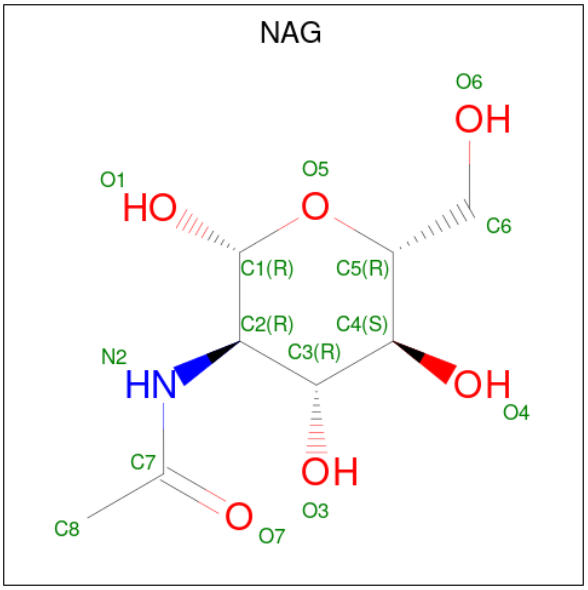
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
			Total	C	N	O			
3	D	3	39	22	2	15	0	0	0
			Total	C	N	O			
3	F	3	39	22	2	15	0	0	0
			Total	C	N	O			
3	H	3	39	22	2	15	0	0	0

- Molecule 4 is an oligosaccharide called alpha-D-mannopyranose-(1-2)-alpha-D-mannopyranose-(1-3)-alpha-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
			Total	C	N	O			
4	G	5	61	34	2	25	0	0	0
			Total	C	N	O			
4	L	5	61	34	2	25	0	0	0

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



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

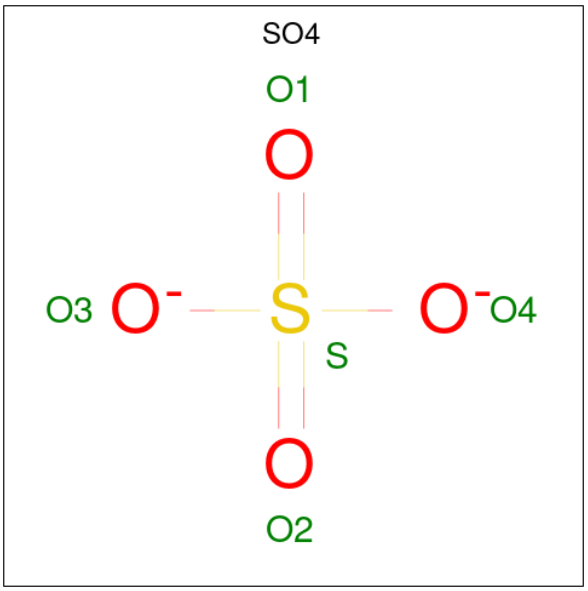
- Molecule 6 is COPPER (II) ION (three-letter code: CU) (formula: Cu).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	Cu		
6	A	4	4	4	0	0
6	B	4	4	4	0	0

- Molecule 7 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

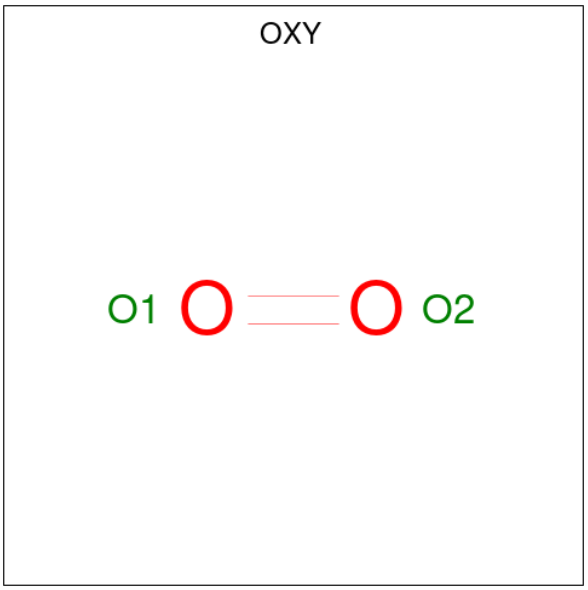
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	Cl		
7	A	1	1	1	0	0
7	B	1	1	1	0	0

- Molecule 8 is SULFATE ION (three-letter code: SO4) (formula: O_4S).



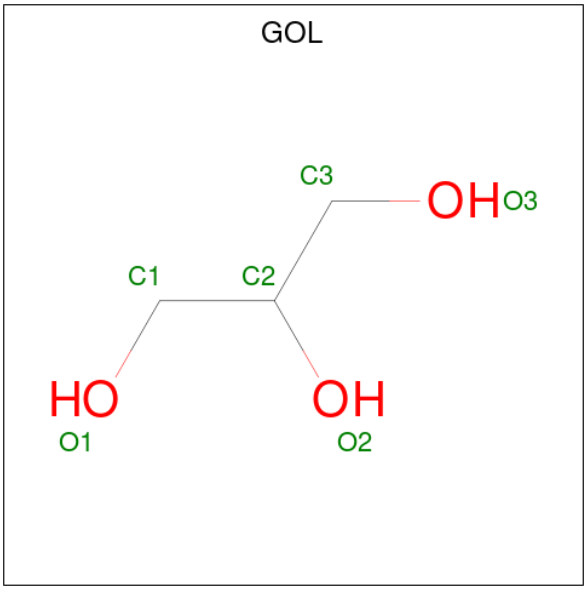
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
8	A	1	Total O S 5 4 1	0	0
8	A	1	Total O S 5 4 1	0	0
8	A	1	Total O S 5 4 1	0	0
8	B	1	Total O S 5 4 1	0	0
8	B	1	Total O S 5 4 1	0	0

- Molecule 9 is OXYGEN MOLECULE (three-letter code: OXY) (formula: O_2).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
9	A	1	Total O 2 2	0	0
9	B	1	Total O 2 2	0	0

- Molecule 10 is GLYCEROL (three-letter code: GOL) (formula: $C_3H_8O_3$).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
10	A	1	Total C O 12 6 6	0	1
10	B	1	Total C O 12 6 6	0	1

- Molecule 11 is water.

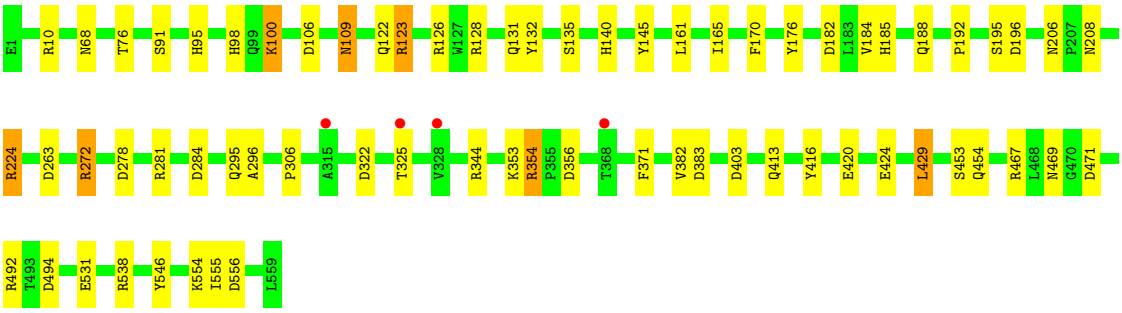
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
11	A	730	Total O 730 730	0	0
11	B	713	Total O 713 713	0	0

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.

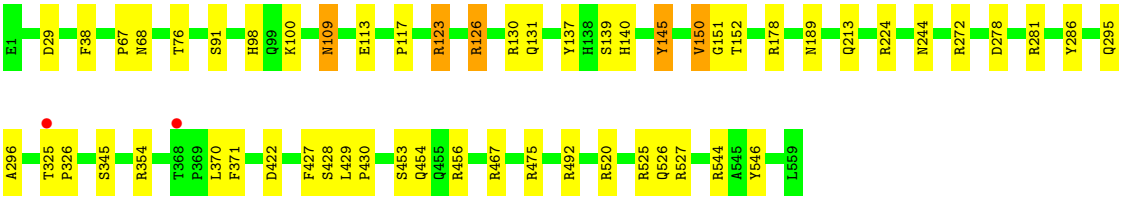
• Molecule 1: Laccase-1

Chain A:



• Molecule 1: Laccase-1

Chain B:

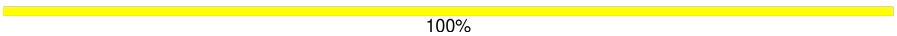


• Molecule 2: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain C:

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

Chain E:



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

Chain I: 100%

NAG1
NAG2

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

Chain J: 100%

NAG1
NAG2

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

Chain K: 100%

NAG1
NAG2

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

Chain D: 100%

NAG1
NAG2
MAN3

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

Chain F: 100%

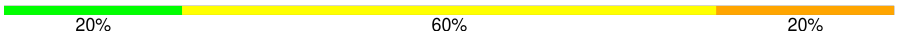
NAG1
NAG2
MAN3

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

Chain H: 100%

NAG1
NAG2
MAN3

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

Chain G: 

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

Chain L:

4 Data and refinement statistics (i)

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α, β, γ	173.74Å 61.90Å 124.21Å 90.00° 96.36° 90.00°	Depositor
Resolution (Å)	15.00 – 1.30 19.63 – 1.30	Depositor EDS
% Data completeness (in resolution range)	91.9 (15.00-1.30) 96.7 (19.63-1.30)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.08	Depositor
$< I/\sigma(I) >$ [1]	1.45 (at 1.30Å)	Xtriage
Refinement program	SHELX, SHELXL-97	Depositor
R, R_{free}	0.167 , 0.212 0.167 , 0.207	Depositor DCC
R_{free} test set	15533 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å^2)	8.3	Xtriage
Anisotropy	0.110	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.34 , 62.3	EDS
L-test for twinning[2]	$< \|L\| > = 0.41$, $< L^2 > = 0.24$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.96	EDS
Total number of atoms	10805	wwPDB-VP
Average B, all atoms (Å^2)	13.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.38% of the height of the origin peak. No significant pseudotranslation is detected.*

[1]Intensities estimated from amplitudes.

[2]Theoretical values of $< |L| >$, $< L^2 >$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality (i)

5.1 Standard geometry (i)

Bond lengths and bond angles in the following residue types are not validated in this section: NAG, OXY, CU, SO4, MAN, GOL, OHI, CL

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	0.55	0/4613	1.28	35/6331 (0.6%)
1	B	0.54	0/4608	1.26	30/6326 (0.5%)
All	All	0.55	0/9221	1.27	65/12657 (0.5%)

There are no bond length outliers.

All (65) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	B	123	ARG	NE-CZ-NH1	-12.91	113.85	120.30
1	A	123	ARG	NE-CZ-NH1	-11.77	114.42	120.30
1	A	128	ARG	NE-CZ-NH2	-11.76	114.42	120.30
1	B	544	ARG	NE-CZ-NH1	10.64	125.62	120.30
1	B	467	ARG	NE-CZ-NH2	-10.62	114.99	120.30
1	B	123	ARG	NE-CZ-NH2	10.54	125.57	120.30
1	A	546	TYR	CB-CG-CD1	-9.72	115.17	121.00
1	B	544	ARG	NE-CZ-NH2	-9.42	115.59	120.30
1	B	281	ARG	CD-NE-CZ	9.00	136.19	123.60
1	B	281	ARG	NE-CZ-NH1	8.56	124.58	120.30
1	B	527	ARG	NE-CZ-NH2	-8.50	116.05	120.30
1	B	546	TYR	CB-CG-CD1	-8.34	115.99	121.00
1	B	467	ARG	NH1-CZ-NH2	8.31	128.54	119.40
1	A	546	TYR	CB-CG-CD2	7.90	125.74	121.00
1	B	286	TYR	CB-CG-CD1	-7.86	116.28	121.00
1	B	525	ARG	NE-CZ-NH2	-7.75	116.42	120.30
1	A	10	ARG	NE-CZ-NH2	-7.66	116.47	120.30
1	B	467	ARG	NE-CZ-NH1	-7.66	116.47	120.30
1	B	520	ARG	NE-CZ-NH2	-7.47	116.57	120.30
1	B	38	PHE	CB-CG-CD2	-7.46	115.58	120.80
1	B	546	TYR	CB-CG-CD2	7.38	125.42	121.00
1	B	492	ARG	NE-CZ-NH2	-7.36	116.62	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	132	TYR	CB-CG-CD2	-7.32	116.61	121.00
1	A	531	GLU	OE1-CD-OE2	-7.28	114.56	123.30
1	A	492	ARG	NE-CZ-NH2	-7.06	116.77	120.30
1	A	145	TYR	CA-CB-CG	6.85	126.41	113.40
1	B	126[A]	ARG	NE-CZ-NH2	6.76	123.68	120.30
1	B	126[B]	ARG	NE-CZ-NH2	6.76	123.68	120.30
1	A	132	TYR	CG-CD2-CE2	-6.76	115.89	121.30
1	A	10	ARG	NE-CZ-NH1	6.48	123.54	120.30
1	B	130	ARG	NE-CZ-NH1	-6.48	117.06	120.30
1	B	354	ARG	NE-CZ-NH1	6.46	123.53	120.30
1	A	224	ARG	NE-CZ-NH1	6.28	123.44	120.30
1	B	145	TYR	CA-CB-CG	6.27	125.31	113.40
1	A	272	ARG	NE-CZ-NH2	-6.16	117.22	120.30
1	A	281	ARG	CD-NE-CZ	6.08	132.12	123.60
1	A	272	ARG	NE-CZ-NH1	6.05	123.33	120.30
1	A	322	ASP	CB-CG-OD2	-5.99	112.91	118.30
1	B	150[A]	VAL	CA-CB-CG2	5.89	119.74	110.90
1	B	150[B]	VAL	CA-CB-CG2	5.89	119.74	110.90
1	A	383	ASP	CB-CG-OD2	-5.89	113.00	118.30
1	B	475	ARG	NE-CZ-NH1	5.86	123.23	120.30
1	A	429[A]	LEU	CA-CB-CG	5.78	128.60	115.30
1	A	429[B]	LEU	CA-CB-CG	5.78	128.60	115.30
1	B	525	ARG	NE-CZ-NH1	5.75	123.18	120.30
1	B	178	ARG	NE-CZ-NH2	5.61	123.11	120.30
1	A	356	ASP	CB-CG-OD2	-5.60	113.26	118.30
1	B	113	GLU	OE1-CD-OE2	-5.57	116.62	123.30
1	A	354	ARG	NE-CZ-NH1	-5.56	117.52	120.30
1	B	422	ASP	CB-CG-OD1	5.48	123.23	118.30
1	A	322	ASP	CB-CG-OD1	5.47	123.23	118.30
1	A	196	ASP	CB-CG-OD2	5.42	123.18	118.30
1	B	286	TYR	CG-CD1-CE1	-5.41	116.97	121.30
1	A	403	ASP	CB-CG-OD2	-5.40	113.44	118.30
1	A	106	ASP	CB-CG-OD1	5.38	123.14	118.30
1	A	284	ASP	CB-CG-OD1	5.37	123.13	118.30
1	A	538	ARG	CD-NE-CZ	5.34	131.07	123.60
1	A	68[A]	ASN	N-CA-CB	5.31	120.16	110.60
1	A	68[B]	ASN	N-CA-CB	5.31	120.16	110.60
1	A	344	ARG	NE-CZ-NH1	5.26	122.93	120.30
1	A	128	ARG	NE-CZ-NH1	5.20	122.90	120.30
1	A	494	ASP	CB-CG-OD2	5.16	122.94	118.30
1	A	416	TYR	CB-CG-CD1	-5.14	117.92	121.00
1	A	263	ASP	CB-CG-OD2	-5.02	113.78	118.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	467	ARG	NE-CZ-NH2	-5.02	117.79	120.30

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	4432	0	4171	37	0
1	B	4418	0	4154	35	0
2	C	28	0	25	4	0
2	E	28	0	25	0	0
2	I	28	0	25	0	0
2	J	28	0	25	0	0
2	K	28	0	25	0	0
3	D	39	0	34	0	0
3	F	39	0	34	0	0
3	H	39	0	34	0	0
4	G	61	0	52	2	0
4	L	61	0	52	0	0
5	A	42	0	39	0	0
5	B	28	0	26	0	0
6	A	4	0	0	0	0
6	B	4	0	0	0	0
7	A	1	0	0	0	0
7	B	1	0	0	0	0
8	A	15	0	0	0	0
8	B	10	0	0	0	0
9	A	2	0	0	0	0
9	B	2	0	0	0	0
10	A	12	0	16	2	0
10	B	12	0	16	3	0
11	A	730	0	0	17	0
11	B	713	0	0	14	0
All	All	10805	0	8753	76	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 4.

All (76) 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:126[A]:ARG:NH1	11:A:1105:HOH:O	1.60	1.34
1:A:126[A]:ARG:CZ	11:A:1105:HOH:O	1.82	1.23
1:A:182[B]:ASP:OD1	11:A:1106:HOH:O	1.76	1.01
2:C:1:NAG:C4	2:C:2:NAG:C1	2.53	0.86
1:A:161:LEU:HG	1:A:224:ARG:HH12	1.41	0.84
1:A:126[A]:ARG:NH2	11:A:1105:HOH:O	1.95	0.80
1:A:192:PRO:HG3	11:A:1694:HOH:O	1.84	0.76
1:A:295:GLN:OE1	1:B:295:GLN:HA	1.86	0.76
1:B:295:GLN:HG2	11:B:1504:HOH:O	1.88	0.74
1:B:126[B]:ARG:NH1	11:B:1105:HOH:O	1.73	0.72
1:B:67:PRO:HG3	11:B:1222:HOH:O	1.89	0.72
2:C:1:NAG:O4	2:C:2:NAG:C2	2.39	0.71
1:A:165:ILE:HA	11:A:1292:HOH:O	1.92	0.69
1:A:131[C]:GLN:HG3	11:A:1101:HOH:O	1.92	0.69
1:B:429:LEU:HD11	10:B:1011[B]:GOL:H32	1.78	0.66
1:B:145:TYR:HB2	1:B:150[B]:VAL:HG13	1.78	0.66
1:B:213:GLN:HB3	11:B:1645:HOH:O	1.97	0.65
1:B:145:TYR:HB2	1:B:150[A]:VAL:HG22	1.79	0.65
1:A:371:PHE:CD2	10:A:1015[A]:GOL:H32	2.34	0.63
1:B:68[B]:ASN:HB3	11:B:1102:HOH:O	1.97	0.63
1:A:224:ARG:HG3	11:A:1316:HOH:O	1.98	0.63
1:B:131[A]:GLN:HG3	11:B:1479:HOH:O	1.99	0.62
1:A:429[B]:LEU:HD11	10:A:1015[B]:GOL:H31	1.82	0.61
1:B:371:PHE:CG	10:B:1011[B]:GOL:H11	2.37	0.59
1:A:554:LYS:HE3	1:A:556:ASP:O	2.01	0.59
1:A:126[B]:ARG:HD3	11:A:1590:HOH:O	2.03	0.58
11:A:1107:HOH:O	4:G:1:NAG:H4	2.04	0.58
1:B:326:PRO:HB3	11:B:1621:HOH:O	2.06	0.56
1:B:76:THR:HG21	1:B:126[B]:ARG:CZ	2.36	0.56
1:A:161:LEU:CG	1:A:224:ARG:HH12	2.15	0.56
1:A:469:ASN:ND2	1:A:471:ASP:H	2.08	0.52
1:A:76:THR:HG21	1:A:126[A]:ARG:CZ	2.39	0.52
1:B:145:TYR:HB2	1:B:150[B]:VAL:CG1	2.37	0.51
1:A:295:GLN:HA	1:B:295:GLN:HE21	1.77	0.50
1:A:354:ARG:HD3	11:A:1706:HOH:O	2.11	0.50
1:A:353:LYS:HD3	11:A:1428:HOH:O	2.10	0.50
1:A:170:PHE:CD1	4:G:1:NAG:H82	2.47	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:224:ARG:HD3	1:A:278[B]:ASP:OD1	2.13	0.48
1:A:100:LYS:HB2	1:A:100:LYS:HE3	1.56	0.48
1:B:526:GLN:HG3	11:B:1179:HOH:O	2.13	0.48
1:A:555:ILE:HG12	11:A:1625:HOH:O	2.14	0.47
1:A:91[B]:SER:OG	1:A:109:ASN:HB3	2.14	0.47
1:A:353:LYS:HE3	1:A:420:GLU:OE2	2.15	0.47
1:B:189:ASN:ND2	1:B:370:LEU:H	2.13	0.46
1:B:456:ARG:NE	11:B:1139:HOH:O	2.49	0.46
1:B:117:PRO:HD2	11:B:1489:HOH:O	2.15	0.45
1:B:91[B]:SER:OG	1:B:109:ASN:HB3	2.17	0.45
1:B:430:PRO:HB3	11:B:1382:HOH:O	2.16	0.45
1:A:122:GLN:NE2	11:A:1153:HOH:O	2.50	0.45
1:A:413[B]:GLN:NE2	11:A:1146:HOH:O	2.49	0.45
1:A:295:GLN:O	1:A:296:ALA:HB3	2.17	0.45
1:B:295:GLN:O	1:B:296:ALA:HB3	2.17	0.45
1:B:137:TYR:CZ	1:B:151:GLY:HA3	2.53	0.44
1:A:382:VAL:O	1:A:554:LYS:NZ	2.49	0.44
1:B:371:PHE:CD2	10:B:1011[B]:GOL:H11	2.52	0.44
1:A:100:LYS:NZ	11:A:1147:HOH:O	2.50	0.44
1:B:68[A]:ASN:ND2	1:B:152:THR:OG1	2.50	0.44
1:B:427:PHE:C	1:B:456:ARG:HH22	2.22	0.43
1:B:428:SER:C	1:B:429:LEU:HD23	2.39	0.43
1:B:29:ASP:O	1:B:29:ASP:OD2	2.38	0.42
1:A:176:TYR:CE2	1:A:195:SER:HA	2.53	0.42
1:B:454:GLN:NE2	11:B:1144:HOH:O	2.50	0.42
1:A:206:ASN:OD1	1:A:208:ASN:HB2	2.19	0.42
1:A:453:SER:O	1:A:454:GLN:HB2	2.18	0.42
1:B:224:ARG:HG2	1:B:278:ASP:OD1	2.18	0.42
1:B:325:THR:HB	1:B:326:PRO:HD2	2.00	0.42
2:C:1:NAG:H61	2:C:2:NAG:H82	2.01	0.42
1:A:184:VAL:O	1:A:188:GLN:HG3	2.19	0.42
1:B:244:ASN:HA	11:B:1109:HOH:O	2.19	0.42
1:A:95:HIS:O	1:A:135:SER:HB3	2.19	0.41
1:B:453:SER:O	1:B:454:GLN:HB2	2.19	0.41
2:C:1:NAG:H4	2:C:2:NAG:C1	2.43	0.41
1:A:185:HIS:HD2	11:A:1336:HOH:O	2.03	0.41
1:B:272:ARG:N	1:B:272:ARG:HD3	2.36	0.40

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	571/559 (102%)	559 (98%)	12 (2%)	0	100	100
1	B	571/559 (102%)	559 (98%)	12 (2%)	0	100	100
All	All	1142/1118 (102%)	1118 (98%)	24 (2%)	0	100	100

There are no Ramachandran outliers to report.

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	492/477 (103%)	484 (98%)	8 (2%)	62	28
1	B	492/477 (103%)	487 (99%)	5 (1%)	76	48
All	All	984/954 (103%)	971 (99%)	13 (1%)	67	35

All (13) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	100	LYS
1	A	109	ASN
1	A	123	ARG
1	A	140	HIS
1	A	272	ARG
1	A	306	PRO
1	A	325	THR
1	A	424	GLU

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Mol	Chain	Res	Type
1	B	100	LYS
1	B	109	ASN
1	B	123	ARG
1	B	140	HIS
1	B	345	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (13) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	34	GLN
1	A	99	GLN
1	A	122	GLN
1	A	185	HIS
1	A	469	ASN
1	B	84	ASN
1	B	99	GLN
1	B	185	HIS
1	B	189	ASN
1	B	295	GLN
1	B	454	GLN
1	B	455	GLN
1	B	526	GLN

5.3.3 RNA (i)

There are no RNA molecules in this entry.

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

2 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	$\#\|Z\| > 2$
1	OHI	B	98	-	9,11,12	1.47	1 (11%)	3,14,16	1.93	2 (66%)
1	OHI	A	98	-	9,11,12	1.42	0	3,14,16	2.30	1 (33%)

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
1	OHI	B	98	-	-	0/4/15/17	0/1/1/1
1	OHI	A	98	-	-	0/4/15/17	0/1/1/1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	98	OHI	O12-CE1	2.22	1.29	1.24

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	98	OHI	O12-CE1-ND1	-3.43	117.14	125.86
1	B	98	OHI	O12-CE1-ND1	-2.41	119.72	125.86
1	B	98	OHI	CB-CG-CD2	-2.30	119.19	129.27

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.5 Carbohydrates (i)

29 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	C	1	2,1	14,14,15	1.15	1 (7%)	17,19,21	1.48	3 (17%)
2	NAG	C	2	2	14,14,15	0.84	1 (7%)	17,19,21	1.49	3 (17%)
3	NAG	D	1	3,1	14,14,15	0.72	0	17,19,21	1.17	1 (5%)
3	NAG	D	2	3	14,14,15	0.87	1 (7%)	17,19,21	1.36	4 (23%)
3	MAN	D	3	3	11,11,12	0.86	0	15,15,17	1.90	6 (40%)
2	NAG	E	1	2,1	14,14,15	0.95	1 (7%)	17,19,21	2.08	3 (17%)
2	NAG	E	2	2	14,14,15	0.92	1 (7%)	17,19,21	1.71	5 (29%)
3	NAG	F	1	3,1	14,14,15	0.83	1 (7%)	17,19,21	1.35	3 (17%)
3	NAG	F	2	3	14,14,15	0.81	1 (7%)	17,19,21	1.47	3 (17%)
3	MAN	F	3	3	11,11,12	0.87	1 (9%)	15,15,17	2.25	5 (33%)
4	NAG	G	1	4	14,14,15	1.16	2 (14%)	17,19,21	2.14	8 (47%)
4	NAG	G	2	4	14,14,15	0.84	1 (7%)	17,19,21	1.09	1 (5%)
4	MAN	G	3	4	11,11,12	0.58	0	15,15,17	0.86	0
4	MAN	G	4	4	11,11,12	0.59	0	15,15,17	1.12	1 (6%)
4	MAN	G	5	4	11,11,12	0.55	0	15,15,17	1.13	1 (6%)
3	NAG	H	1	3,1	14,14,15	0.76	1 (7%)	17,19,21	1.44	1 (5%)
3	NAG	H	2	3	14,14,15	0.73	0	17,19,21	1.58	2 (11%)
3	MAN	H	3	3	11,11,12	0.57	0	15,15,17	1.65	2 (13%)
2	NAG	I	1	2,1	14,14,15	0.76	1 (7%)	17,19,21	1.53	3 (17%)
2	NAG	I	2	2	14,14,15	0.77	0	17,19,21	1.52	4 (23%)
2	NAG	J	1	2,1	14,14,15	0.84	1 (7%)	17,19,21	1.28	1 (5%)
2	NAG	J	2	2	14,14,15	0.77	1 (7%)	17,19,21	1.84	2 (11%)
2	NAG	K	1	2,1	14,14,15	1.21	1 (7%)	17,19,21	1.74	3 (17%)
2	NAG	K	2	2	14,14,15	0.88	1 (7%)	17,19,21	2.38	7 (41%)
4	NAG	L	1	4,1	14,14,15	1.06	1 (7%)	17,19,21	1.34	2 (11%)
4	NAG	L	2	4	14,14,15	0.79	0	17,19,21	1.35	1 (5%)
4	MAN	L	3	4	11,11,12	0.63	0	15,15,17	0.87	0
4	MAN	L	4	4	11,11,12	0.97	1 (9%)	15,15,17	1.57	3 (20%)
4	MAN	L	5	4	11,11,12	0.73	0	15,15,17	1.53	1 (6%)

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

All (19) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	K	1	NAG	C1-C2	3.98	1.57	1.52
2	C	1	NAG	C1-C2	3.69	1.57	1.52
4	G	1	NAG	C1-C2	3.49	1.57	1.52
4	L	1	NAG	C1-C2	2.92	1.56	1.52
2	E	1	NAG	C1-C2	2.65	1.56	1.52
4	G	2	NAG	C1-C2	2.63	1.55	1.52
2	C	2	NAG	C1-C2	2.58	1.55	1.52
2	J	1	NAG	C1-C2	2.58	1.55	1.52
2	E	2	NAG	C1-C2	2.57	1.55	1.52

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	D	2	NAG	C1-C2	2.53	1.55	1.52
4	L	4	MAN	C2-C3	2.43	1.56	1.52
2	K	2	NAG	C1-C2	2.43	1.55	1.52
3	F	2	NAG	C1-C2	2.29	1.55	1.52
2	J	2	NAG	C1-C2	2.28	1.55	1.52
2	I	1	NAG	C1-C2	2.20	1.55	1.52
3	H	1	NAG	C1-C2	2.16	1.55	1.52
3	F	3	MAN	O5-C1	-2.13	1.40	1.43
3	F	1	NAG	C1-C2	2.08	1.55	1.52
4	G	1	NAG	C3-C2	2.01	1.56	1.52

All (79) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	K	2	NAG	C2-N2-C7	-5.56	115.45	122.90
2	K	1	NAG	O5-C1-C2	-5.33	103.04	111.29
2	E	1	NAG	C2-N2-C7	-4.99	116.21	122.90
3	F	3	MAN	C1-O5-C5	4.77	118.57	112.19
4	L	5	MAN	C1-O5-C5	4.61	118.36	112.19
2	I	1	NAG	O5-C1-C2	-4.54	104.27	111.29
2	K	2	NAG	O7-C7-N2	-4.50	114.03	121.98
2	J	2	NAG	O7-C7-C8	-4.47	114.09	122.05
2	E	1	NAG	O5-C1-C2	-4.40	104.48	111.29
3	H	1	NAG	O5-C1-C2	-4.39	104.50	111.29
3	H	3	MAN	O3-C3-C2	4.22	118.66	110.05
2	J	2	NAG	O5-C1-C2	-4.18	104.82	111.29
3	H	2	NAG	O7-C7-C8	-3.91	115.09	122.05
4	G	1	NAG	C1-O5-C5	-3.90	106.95	112.19
2	C	1	NAG	O5-C1-C2	-3.89	105.27	111.29
4	L	1	NAG	O5-C1-C2	-3.77	105.46	111.29
3	F	1	NAG	O5-C1-C2	-3.68	105.59	111.29
3	F	3	MAN	O3-C3-C2	3.61	117.42	110.05
3	D	1	NAG	O5-C1-C2	-3.54	105.82	111.29
2	J	1	NAG	O5-C1-C2	-3.52	105.84	111.29
2	E	1	NAG	O5-C5-C6	3.49	114.46	107.66
2	I	2	NAG	O4-C4-C5	3.43	117.76	109.32
3	D	3	MAN	O5-C5-C6	3.36	114.21	107.66
4	G	1	NAG	O4-C4-C5	-3.35	101.07	109.32
2	E	2	NAG	O4-C4-C3	-3.32	102.54	110.38
4	L	4	MAN	O4-C4-C5	3.29	117.42	109.32
3	F	2	NAG	O5-C1-C2	-3.26	106.25	111.29
3	F	3	MAN	O2-C2-C3	3.18	116.73	110.15

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
4	G	2	NAG	O5-C1-C2	-3.16	106.41	111.29
2	K	2	NAG	C8-C7-N2	3.13	121.31	116.12
2	K	2	NAG	C4-C3-C2	2.98	115.38	111.02
3	D	3	MAN	C2-C3-C4	-2.98	105.62	110.86
4	L	2	NAG	C1-O5-C5	2.93	116.12	112.19
4	G	1	NAG	O5-C1-C2	-2.87	106.86	111.29
2	E	2	NAG	O7-C7-C8	-2.86	116.95	122.05
2	E	2	NAG	C1-O5-C5	2.86	116.02	112.19
4	G	1	NAG	O5-C5-C6	2.85	113.21	107.66
2	I	2	NAG	O7-C7-C8	-2.84	116.99	122.05
2	C	2	NAG	O3-C3-C2	2.84	115.30	109.40
3	H	3	MAN	O5-C5-C6	2.76	113.03	107.66
4	L	4	MAN	C2-C3-C4	-2.76	106.01	110.86
3	F	3	MAN	O4-C4-C3	-2.76	103.88	110.38
2	K	2	NAG	O3-C3-C2	2.75	115.11	109.40
3	D	2	NAG	O5-C1-C2	-2.69	107.14	111.29
3	F	3	MAN	C1-C2-C3	2.67	113.53	109.64
3	F	2	NAG	O5-C5-C4	-2.65	104.39	110.83
2	K	2	NAG	O5-C5-C6	2.62	112.75	107.66
4	L	4	MAN	O2-C2-C3	-2.61	104.74	110.15
2	I	2	NAG	C3-C4-C5	2.61	114.96	110.23
3	D	3	MAN	O4-C4-C3	-2.61	104.23	110.38
2	C	2	NAG	O5-C1-C2	-2.59	107.28	111.29
2	I	2	NAG	O5-C1-C2	-2.54	107.35	111.29
4	G	1	NAG	O7-C7-C8	2.51	126.53	122.05
4	G	5	MAN	C1-C2-C3	2.50	113.28	109.64
3	F	1	NAG	O5-C5-C4	-2.48	104.80	110.83
2	E	2	NAG	C2-N2-C7	-2.45	119.62	122.90
2	E	2	NAG	C4-C3-C2	-2.40	107.51	111.02
4	G	4	MAN	O5-C1-C2	-2.37	105.13	110.79
4	G	1	NAG	O3-C3-C2	-2.35	104.53	109.40
2	C	1	NAG	O5-C5-C4	-2.32	105.19	110.83
3	D	2	NAG	O4-C4-C3	-2.31	104.94	110.38
3	D	3	MAN	O2-C2-C3	-2.25	105.50	110.15
2	K	1	NAG	C2-N2-C7	-2.25	119.89	122.90
2	I	1	NAG	O7-C7-C8	-2.24	118.06	122.05
3	D	2	NAG	C3-C4-C5	-2.24	106.17	110.23
2	C	2	NAG	O3-C3-C4	-2.24	105.10	110.38
2	I	1	NAG	C2-N2-C7	-2.24	119.90	122.90
3	F	2	NAG	C1-O5-C5	-2.23	109.19	112.19
3	D	3	MAN	C1-O5-C5	2.23	115.17	112.19
3	H	2	NAG	C8-C7-N2	2.22	119.80	116.12

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	K	1	NAG	O5-C5-C4	-2.21	105.46	110.83
2	C	1	NAG	C2-N2-C7	2.19	125.84	122.90
4	G	1	NAG	C4-C3-C2	-2.16	107.85	111.02
4	L	1	NAG	O4-C4-C5	-2.16	104.00	109.32
3	F	1	NAG	O3-C3-C2	-2.13	104.98	109.40
3	D	2	NAG	O5-C5-C4	-2.06	105.81	110.83
2	K	2	NAG	C1-C2-N2	-2.06	107.19	110.43
4	G	1	NAG	O7-C7-N2	-2.06	118.35	121.98
3	D	3	MAN	C1-C2-C3	-2.06	106.65	109.64

All (5) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
3	D	3	MAN	C1
3	F	3	MAN	C1
3	H	3	MAN	C1
4	G	3	MAN	C1
4	L	3	MAN	C1

All (17) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	G	5	MAN	O5-C5-C6-O6
3	D	3	MAN	O5-C5-C6-O6
2	K	2	NAG	O5-C5-C6-O6
3	H	3	MAN	O5-C5-C6-O6
3	D	3	MAN	C4-C5-C6-O6
4	G	5	MAN	C4-C5-C6-O6
3	H	3	MAN	C4-C5-C6-O6
2	J	2	NAG	O7-C7-N2-C2
4	L	2	NAG	O7-C7-N2-C2
2	K	2	NAG	C4-C5-C6-O6
2	I	2	NAG	O7-C7-N2-C2
2	C	2	NAG	C8-C7-N2-C2
2	C	2	NAG	O7-C7-N2-C2
2	E	2	NAG	O7-C7-N2-C2
4	L	5	MAN	C4-C5-C6-O6
4	L	4	MAN	O5-C5-C6-O6
4	L	1	NAG	C8-C7-N2-C2

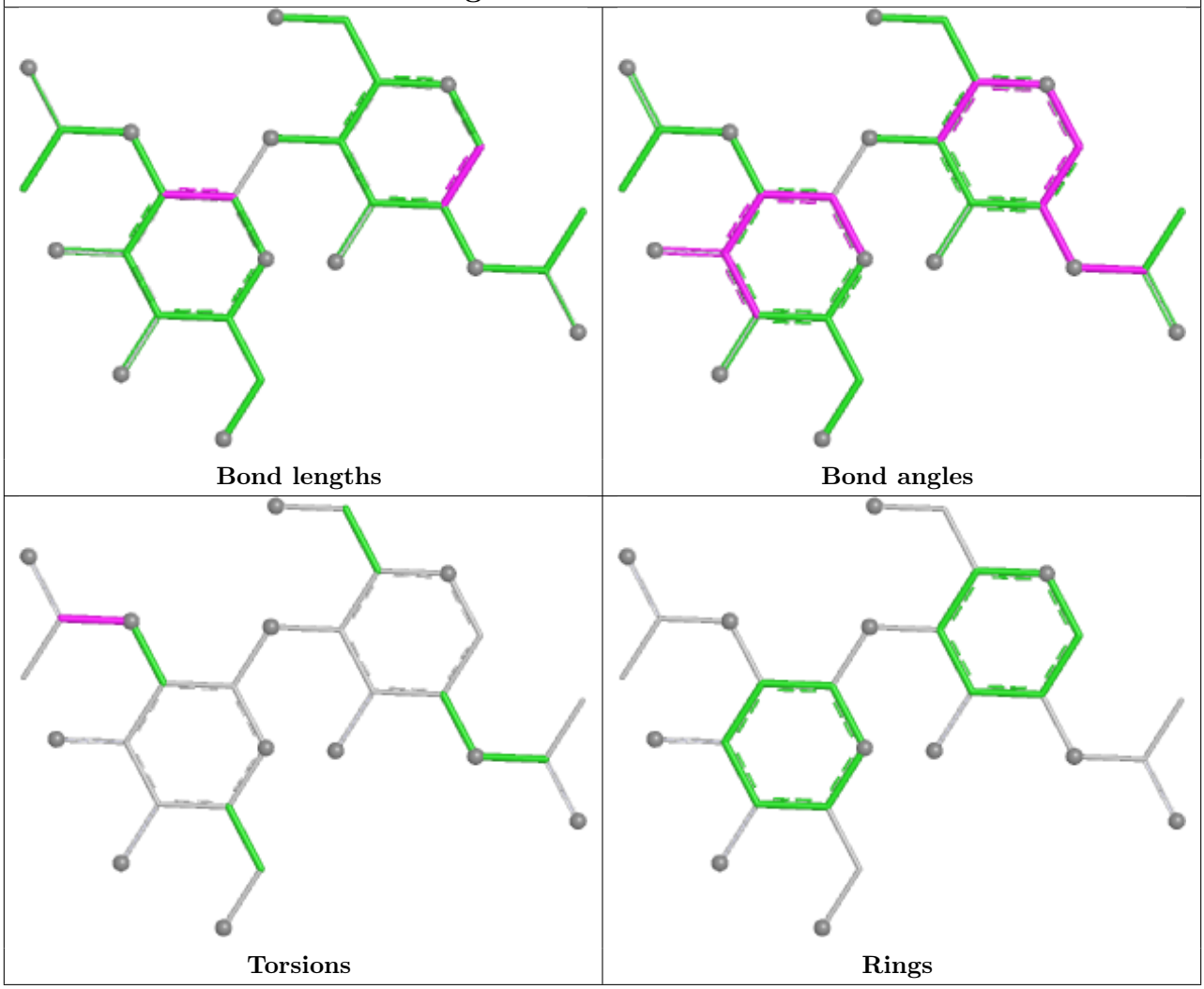
There are no ring outliers.

3 monomers are involved in 6 short contacts:

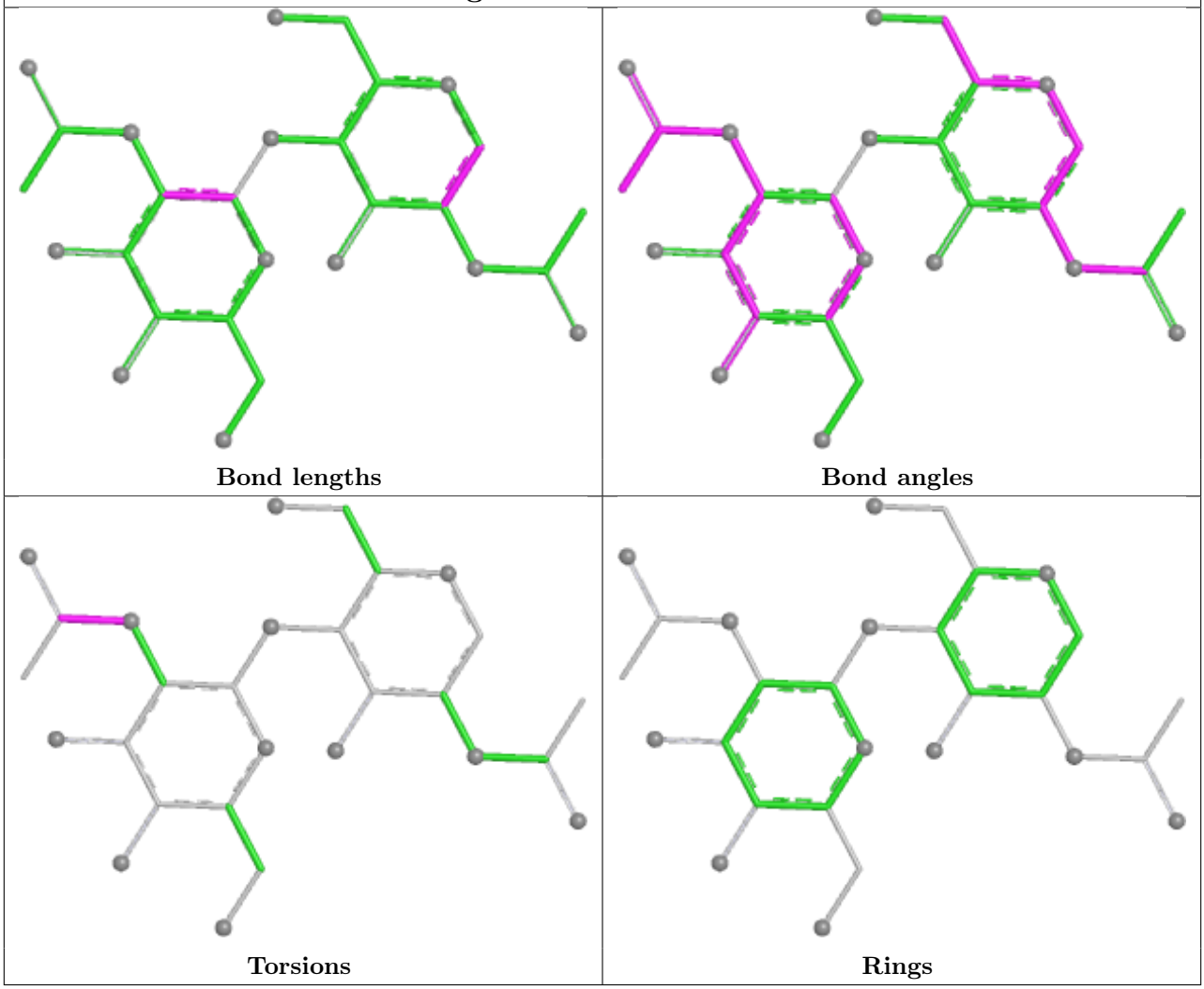
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	2	NAG	4	0
2	C	1	NAG	4	0
4	G	1	NAG	2	0

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

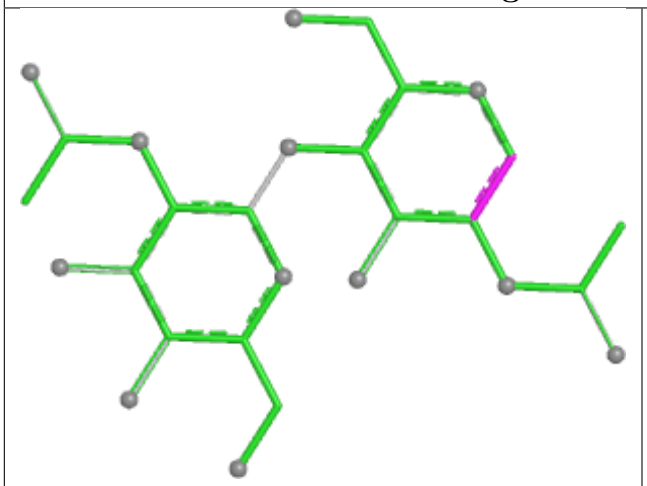
Oligosaccharide Chain C



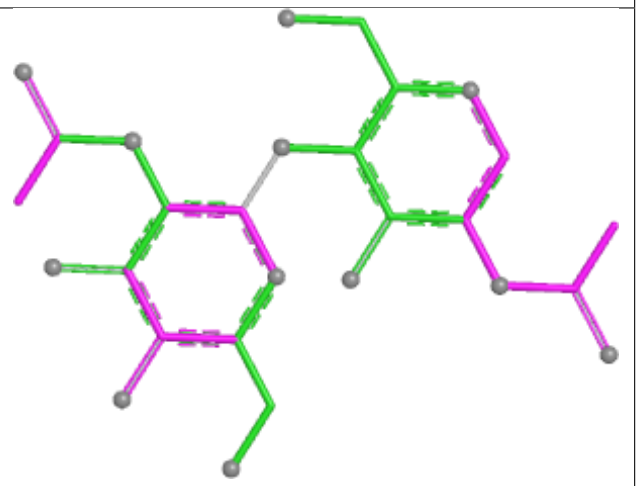
Oligosaccharide Chain E



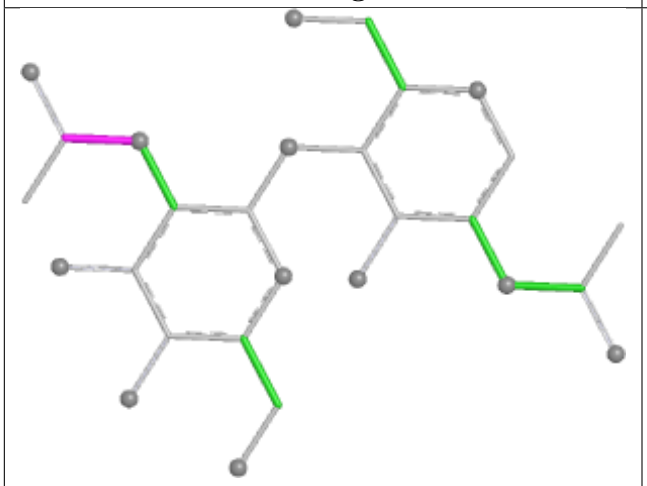
Oligosaccharide Chain I



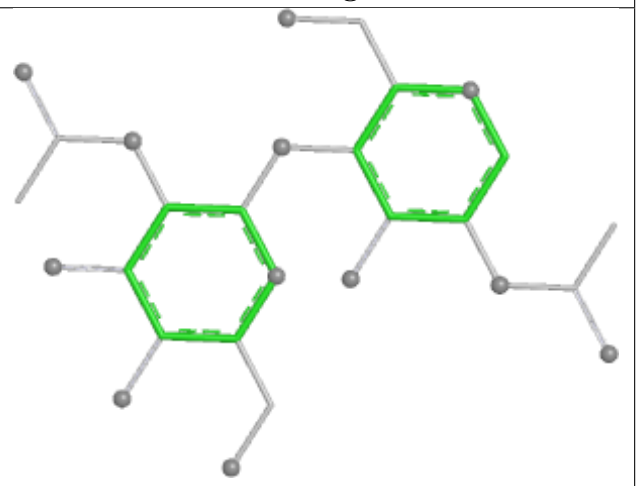
Bond lengths



Bond angles

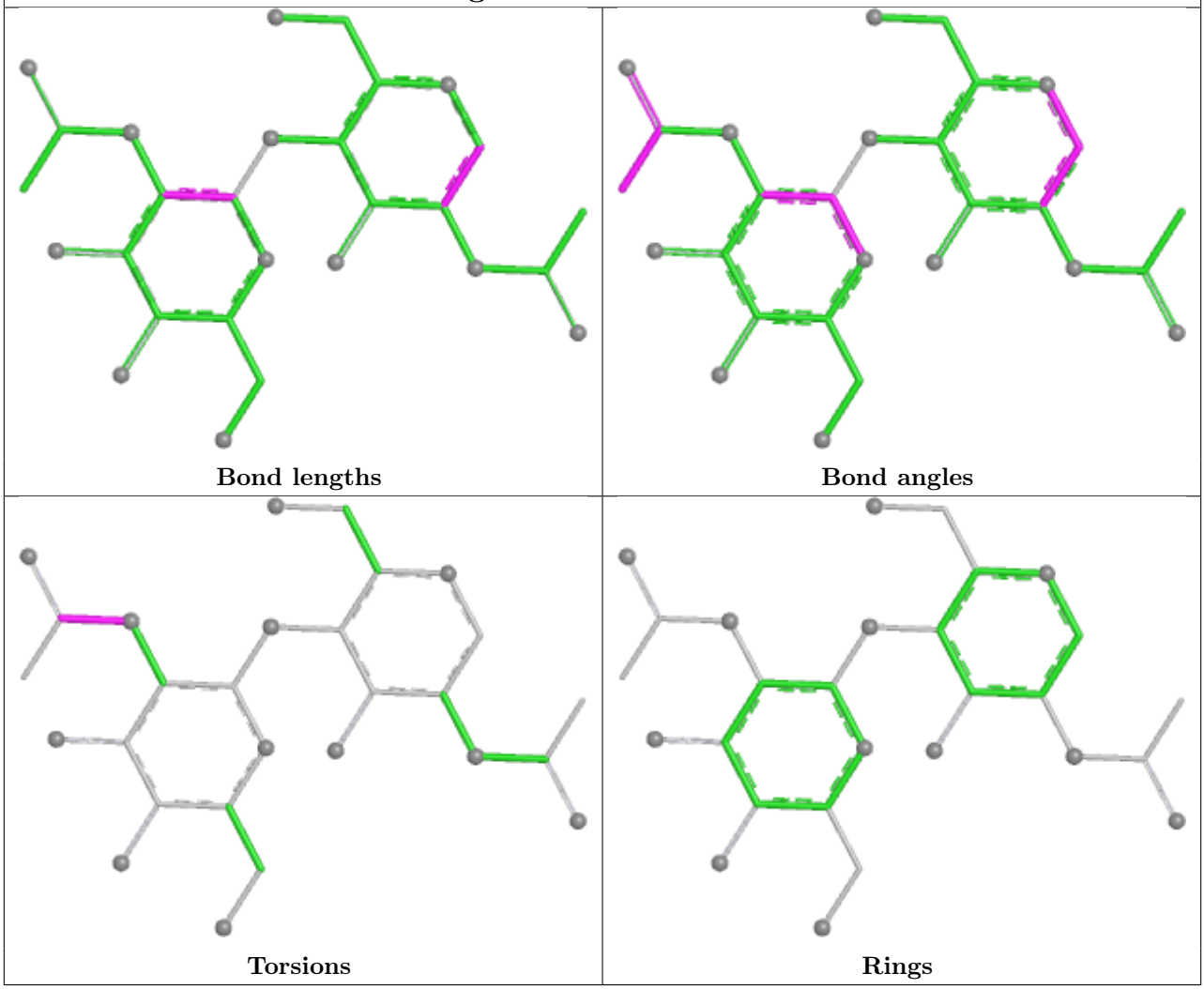


Torsions

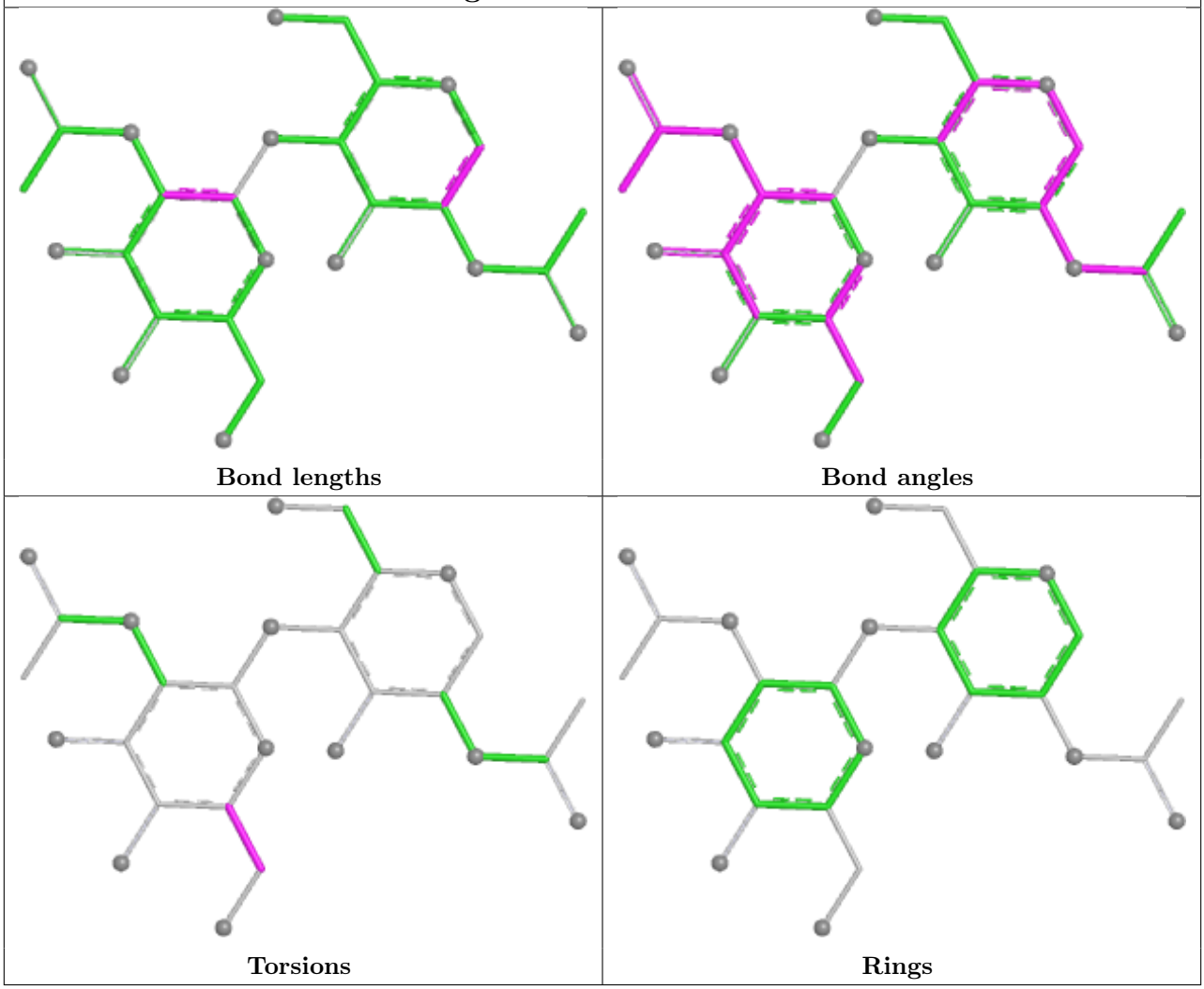


Rings

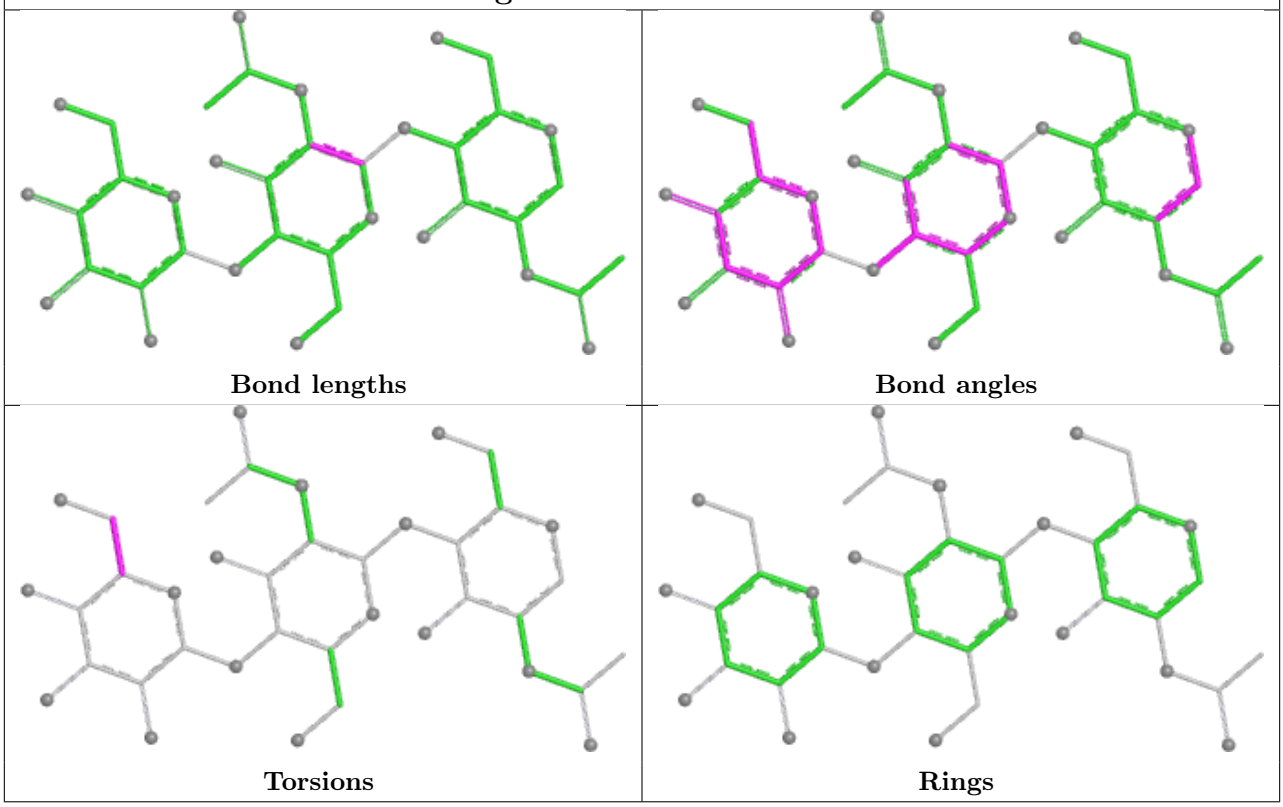
Oligosaccharide Chain J



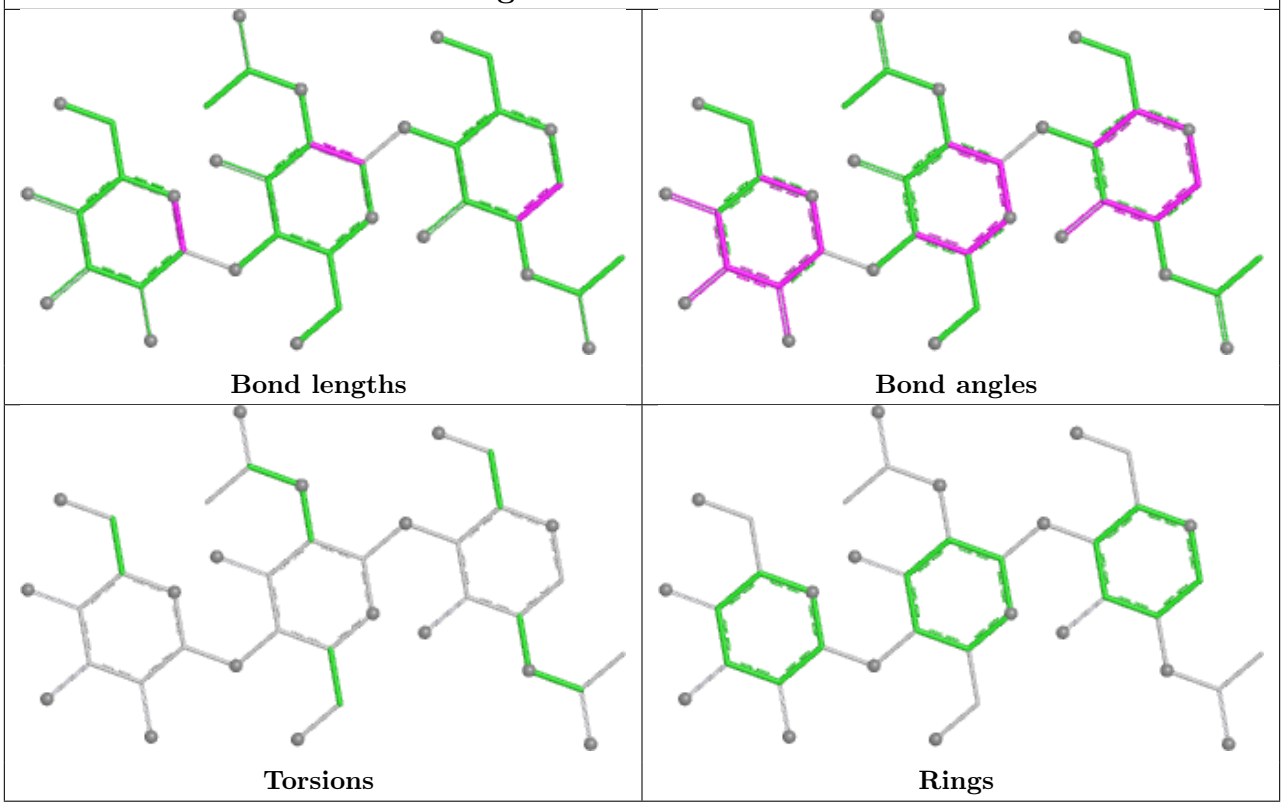
Oligosaccharide Chain K



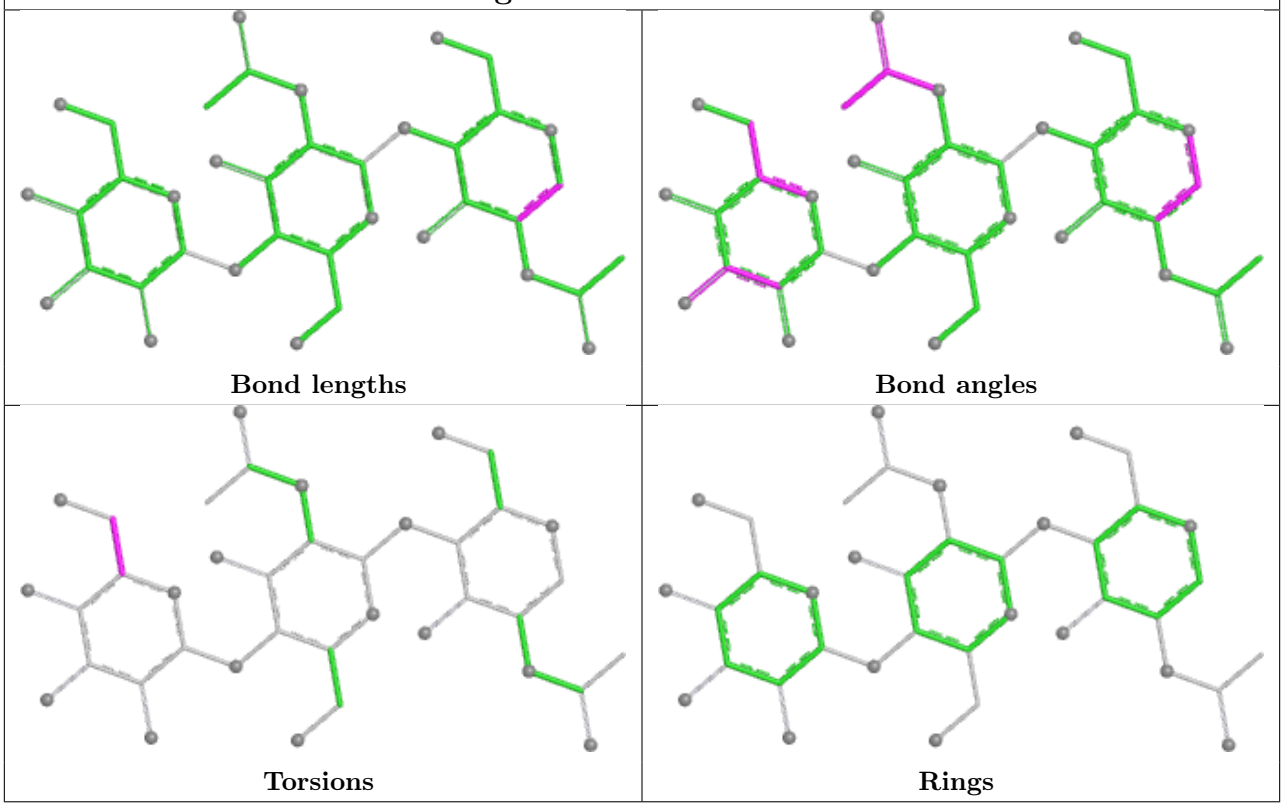
Oligosaccharide Chain D



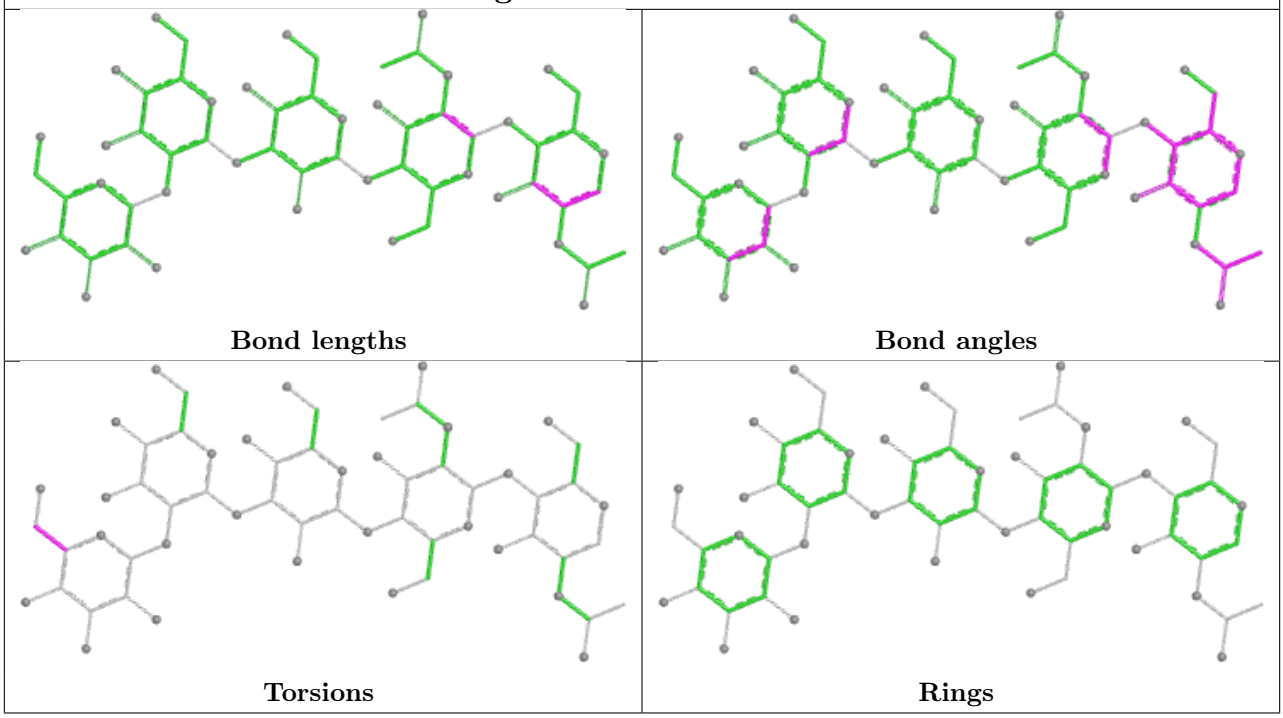
Oligosaccharide Chain F



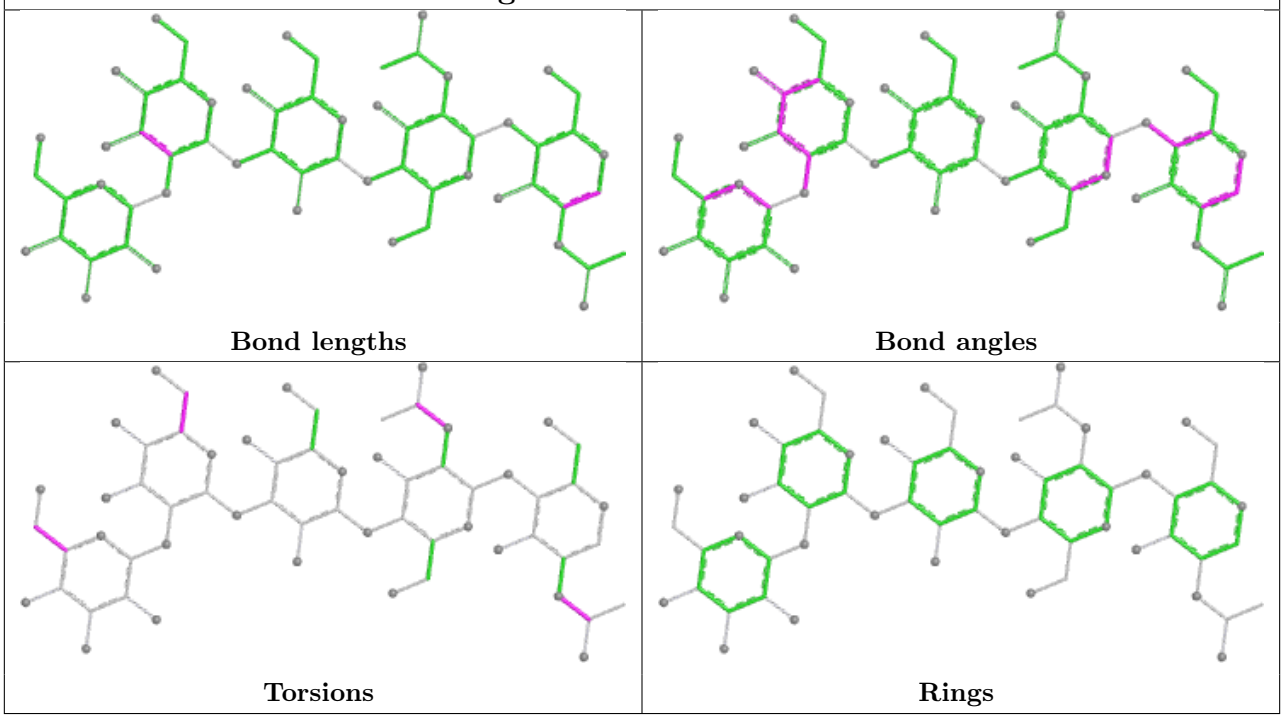
Oligosaccharide Chain H



Oligosaccharide Chain G



Oligosaccharide Chain L



5.6 Ligand geometry (i)

Of 26 ligands modelled in this entry, 10 are monoatomic - leaving 16 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	$\#\|Z\| > 2$
8	SO4	A	1012	-	4,4,4	0.24	0	6,6,6	0.33	0
8	SO4	B	1009	-	4,4,4	0.24	0	6,6,6	0.54	0
8	SO4	A	1011	-	4,4,4	0.18	0	6,6,6	0.28	0
10	GOL	A	1015[B]	-	5,5,5	0.33	0	5,5,5	0.35	0
10	GOL	B	1011[B]	-	5,5,5	0.31	0	5,5,5	0.18	0
10	GOL	A	1015[A]	-	5,5,5	0.25	0	5,5,5	0.89	0
5	NAG	B	1002	1	14,14,15	0.88	1 (7%)	17,19,21	1.26	2 (11%)
5	NAG	B	1001	1	14,14,15	0.81	0	17,19,21	1.58	3 (17%)
9	OXY	A	1014	-	1,1,1	0.66	0	-		
9	OXY	B	1010	-	1,1,1	0.68	0	-		
8	SO4	B	1008	-	4,4,4	0.27	0	6,6,6	0.54	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\| > 2$	Counts	RMSZ	$\#\|Z\| > 2$
5	NAG	A	1004	1	14,14,15	0.60	0	17,19,21	1.70	4 (23%)
8	SO4	A	1013	-	4,4,4	0.21	0	6,6,6	0.25	0
5	NAG	A	1005	1	14,14,15	0.91	1 (7%)	17,19,21	1.36	2 (11%)
5	NAG	A	1001	1	14,14,15	1.05	1 (7%)	17,19,21	2.33	8 (47%)
10	GOL	B	1011[A]	-	5,5,5	0.31	0	5,5,5	0.13	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
10	GOL	B	1011[B]	-	-	2/4/4/4	-
10	GOL	A	1015[B]	-	-	2/4/4/4	-
10	GOL	A	1015[A]	-	-	1/4/4/4	-
5	NAG	B	1002	1	-	0/6/23/26	0/1/1/1
5	NAG	B	1001	1	-	0/6/23/26	0/1/1/1
5	NAG	A	1004	1	-	1/6/23/26	0/1/1/1
5	NAG	A	1005	1	-	1/6/23/26	0/1/1/1
5	NAG	A	1001	1	-	1/6/23/26	0/1/1/1
10	GOL	B	1011[A]	-	-	2/4/4/4	-

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	A	1001	NAG	C1-C2	3.11	1.56	1.52
5	B	1002	NAG	C1-C2	3.00	1.56	1.52
5	A	1005	NAG	C1-C2	2.20	1.55	1.52

All (19) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
5	A	1001	NAG	C2-N2-C7	-5.81	115.11	122.90
5	A	1004	NAG	O7-C7-C8	-4.43	114.17	122.05
5	A	1001	NAG	O5-C5-C6	4.02	115.48	107.66
5	B	1001	NAG	C2-N2-C7	-3.94	117.62	122.90
5	B	1002	NAG	O5-C1-C2	-3.84	105.35	111.29
5	A	1005	NAG	C1-O5-C5	3.37	116.70	112.19
5	A	1004	NAG	O5-C1-C2	-3.09	106.51	111.29
5	A	1001	NAG	O7-C7-N2	-2.72	117.17	121.98
5	A	1001	NAG	O5-C1-C2	-2.72	107.08	111.29

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
5	B	1001	NAG	O5-C5-C4	-2.61	104.47	110.83
5	A	1001	NAG	C1-O5-C5	2.55	115.61	112.19
5	A	1001	NAG	C8-C7-N2	2.52	120.30	116.12
5	B	1002	NAG	O5-C5-C6	2.39	112.32	107.66
5	A	1004	NAG	O5-C5-C6	2.34	112.21	107.66
5	A	1001	NAG	O4-C4-C3	-2.27	105.03	110.38
5	A	1005	NAG	O7-C7-C8	-2.18	118.17	122.05
5	B	1001	NAG	C3-C4-C5	-2.18	106.28	110.23
5	A	1004	NAG	C1-C2-N2	2.11	113.76	110.43
5	A	1001	NAG	O3-C3-C4	-2.02	105.62	110.38

There are no chirality outliers.

All (10) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
10	A	1015[B]	GOL	C1-C2-C3-O3
10	B	1011[A]	GOL	C1-C2-C3-O3
10	B	1011[B]	GOL	O1-C1-C2-C3
10	A	1015[B]	GOL	O2-C2-C3-O3
10	B	1011[A]	GOL	O2-C2-C3-O3
10	B	1011[B]	GOL	O1-C1-C2-O2
10	A	1015[A]	GOL	O1-C1-C2-O2
5	A	1001	NAG	C8-C7-N2-C2
5	A	1004	NAG	C8-C7-N2-C2
5	A	1005	NAG	C8-C7-N2-C2

There are no ring outliers.

3 monomers are involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
10	A	1015[B]	GOL	1	0
10	B	1011[B]	GOL	3	0
10	A	1015[A]	GOL	1	0

5.7 Other polymers (i)

There are no such residues in this entry.

5.8 Polymer linkage issues (i)

There are no chain breaks in this entry.

6 Fit of model and data (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	558/559 (99%)	-0.33	4 (0%)	87	87	6, 10, 20, 28	0
1	B	558/559 (99%)	-0.33	2 (0%)	92	91	6, 11, 20, 26	0
All	All	1116/1118 (99%)	-0.33	6 (0%)	91	91	6, 11, 20, 28	0

All (6) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	368	THR	3.2
1	A	368	THR	2.5
1	A	325	THR	2.3
1	B	325	THR	2.3
1	A	315	ALA	2.1
1	A	328	VAL	2.0

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
1	OHI	B	98	11/12	0.94	0.11	9,12,16,22	0
1	OHI	A	98	11/12	0.95	0.09	8,13,17,23	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

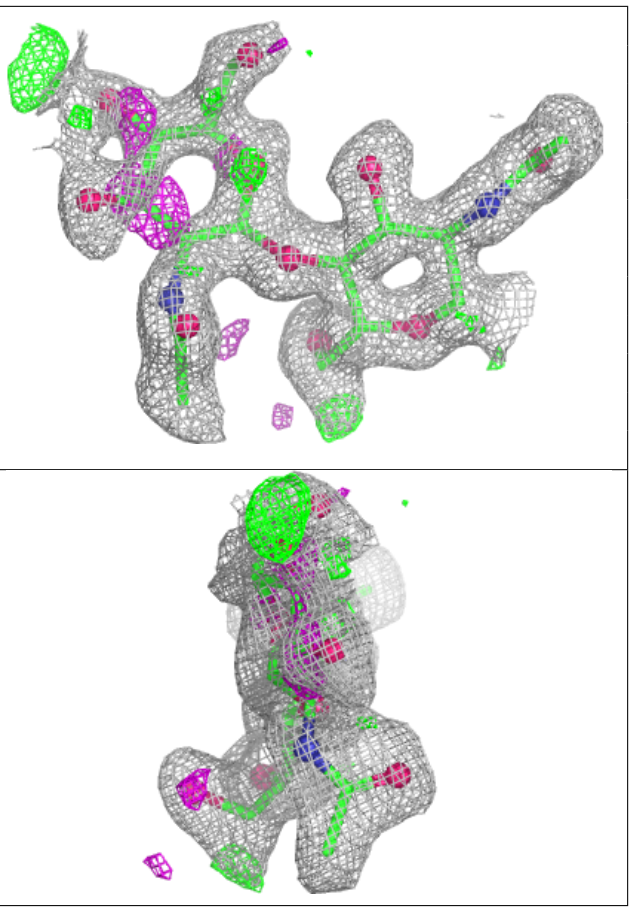
labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
3	MAN	F	3	11/12	0.66	0.33	31,34,37,38	0
2	NAG	C	2	14/15	0.71	0.28	25,28,33,34	0
4	NAG	G	2	14/15	0.73	0.24	23,28,32,32	0
4	MAN	L	5	11/12	0.73	0.17	31,32,33,34	0
4	MAN	G	4	11/12	0.74	0.28	33,34,36,37	0
4	MAN	L	4	11/12	0.75	0.27	33,35,37,37	0
4	NAG	L	2	14/15	0.75	0.21	26,29,30,30	0
4	MAN	G	5	11/12	0.76	0.24	31,32,34,38	0
2	NAG	K	2	14/15	0.81	0.28	25,28,33,34	0
3	MAN	D	3	11/12	0.83	0.16	19,23,27,28	0
3	MAN	H	3	11/12	0.84	0.24	23,26,30,30	0
4	NAG	L	1	14/15	0.85	0.20	19,21,26,31	0
4	NAG	G	1	14/15	0.85	0.19	20,21,26,33	0
4	MAN	L	3	11/12	0.86	0.21	29,30,31,33	0
2	NAG	I	2	14/15	0.86	0.21	24,27,30,32	0
4	MAN	G	3	11/12	0.86	0.22	30,31,32,32	0
3	NAG	F	2	14/15	0.87	0.13	18,21,26,29	0
2	NAG	J	2	14/15	0.90	0.14	17,21,24,25	0
2	NAG	E	2	14/15	0.90	0.17	20,24,26,29	0
2	NAG	I	1	14/15	0.91	0.14	21,23,25,26	0
2	NAG	E	1	14/15	0.91	0.13	17,20,23,26	0
2	NAG	K	1	14/15	0.94	0.08	12,16,22,23	0
3	NAG	H	2	14/15	0.95	0.07	11,14,19,21	0
2	NAG	C	1	14/15	0.96	0.06	12,14,21,23	0
3	NAG	F	1	14/15	0.97	0.06	12,14,16,16	0
3	NAG	D	2	14/15	0.97	0.06	11,13,16,19	0
2	NAG	J	1	14/15	0.97	0.06	13,14,18,18	0
3	NAG	D	1	14/15	0.98	0.06	8,8,11,12	0
3	NAG	H	1	14/15	0.98	0.06	8,9,11,11	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.

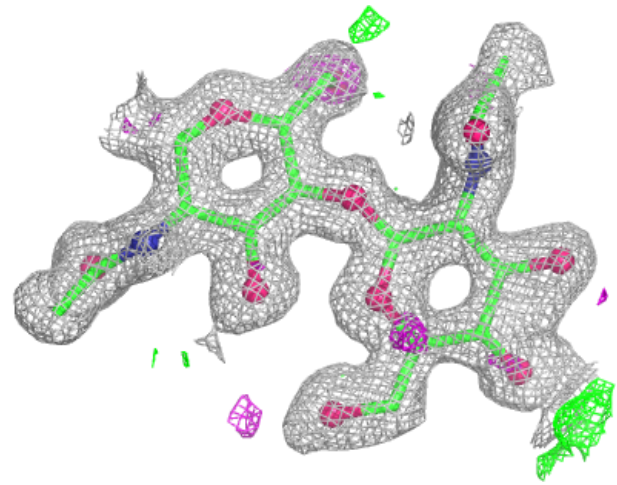
Electron density around Chain C:

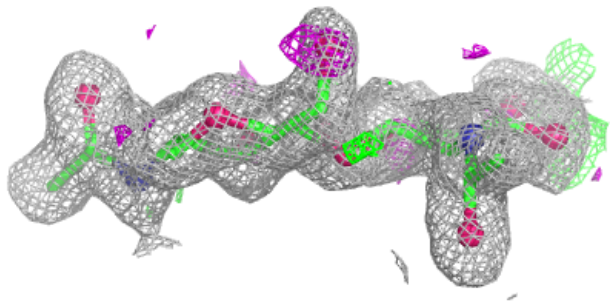
$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

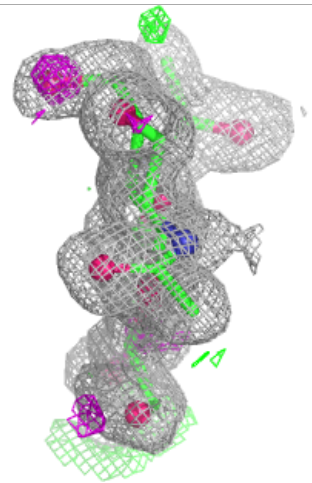


Electron density around Chain E:

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

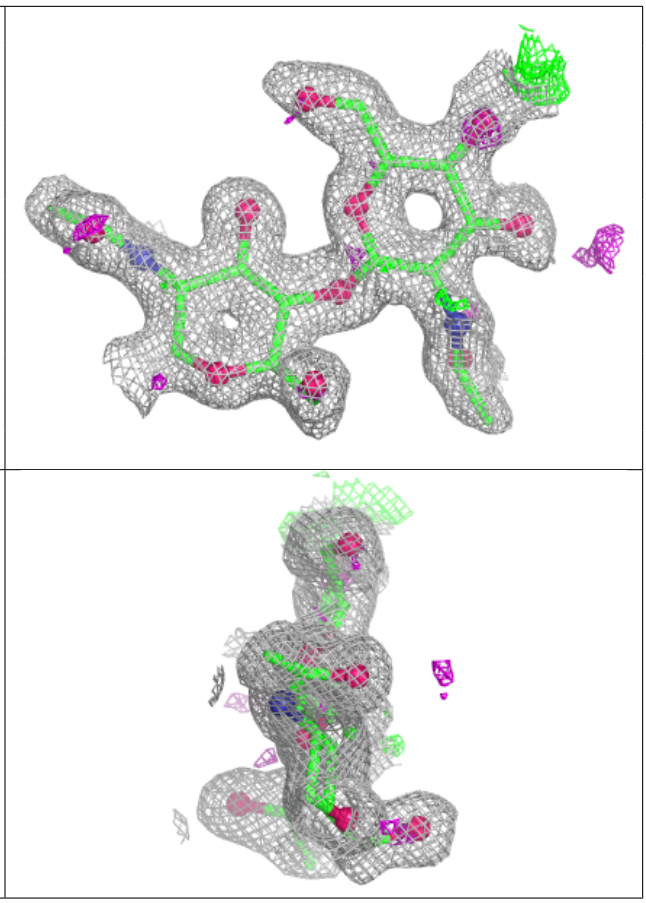






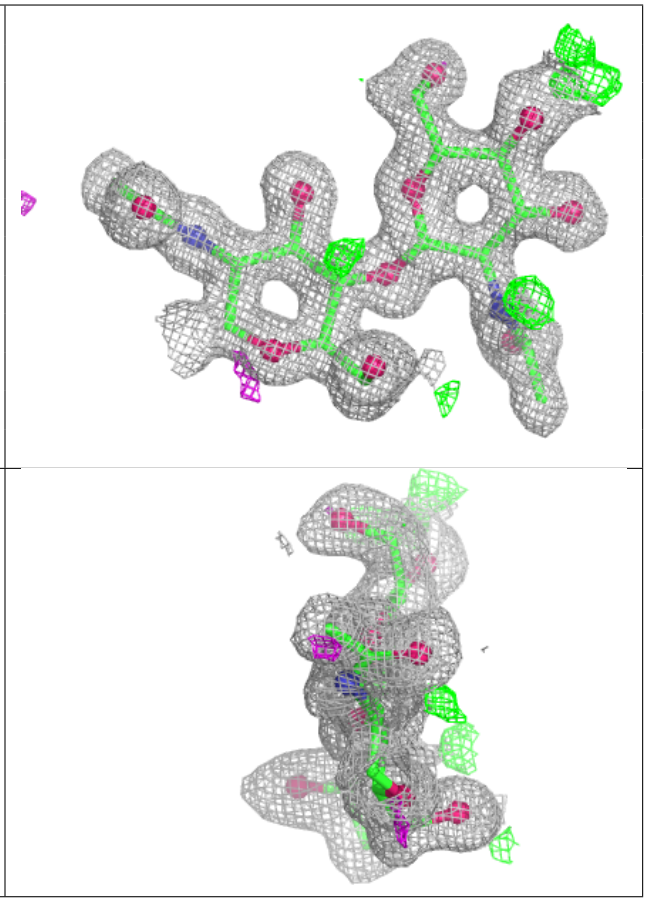
Electron density around Chain I:

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



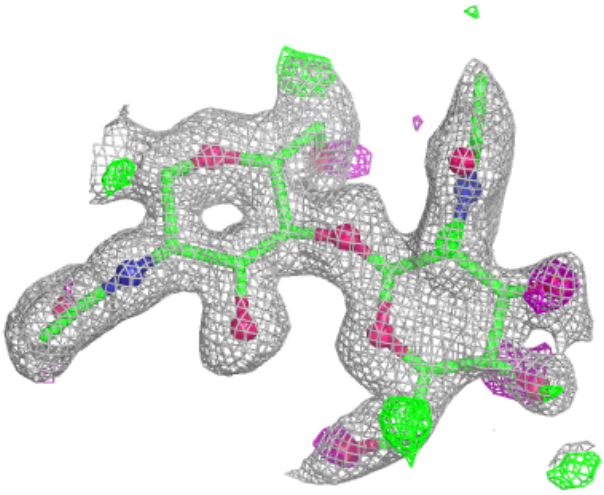
Electron density around Chain J:

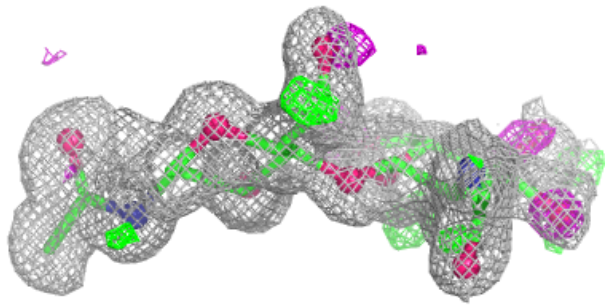
$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

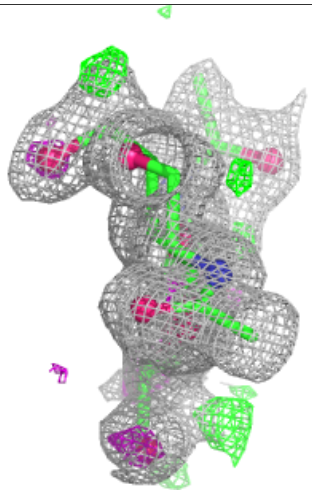


Electron density around Chain K:

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

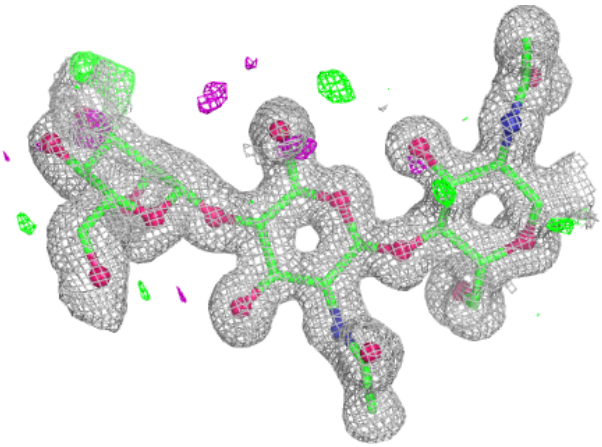


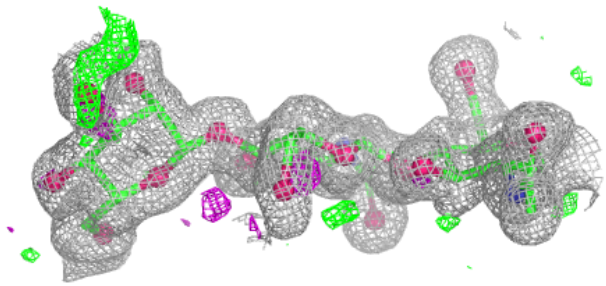


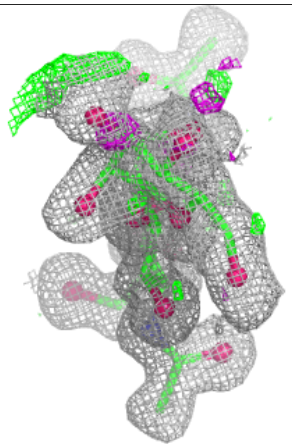


Electron density around Chain D:

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

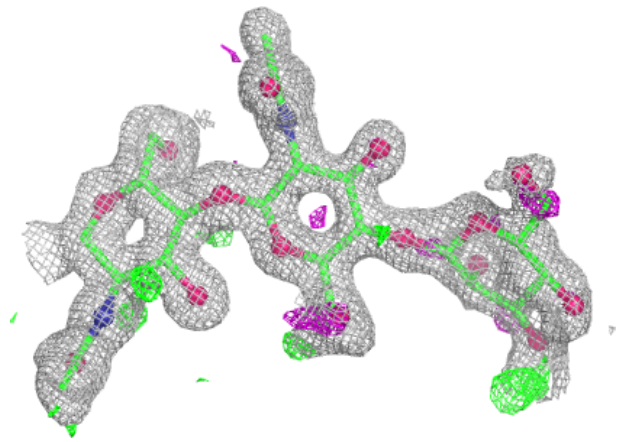


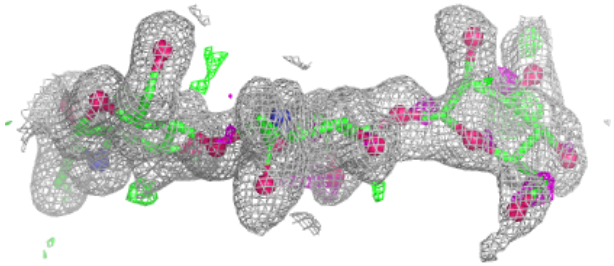


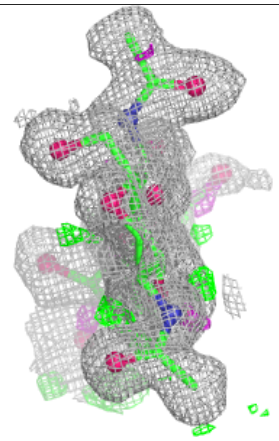


Electron density around Chain F:

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

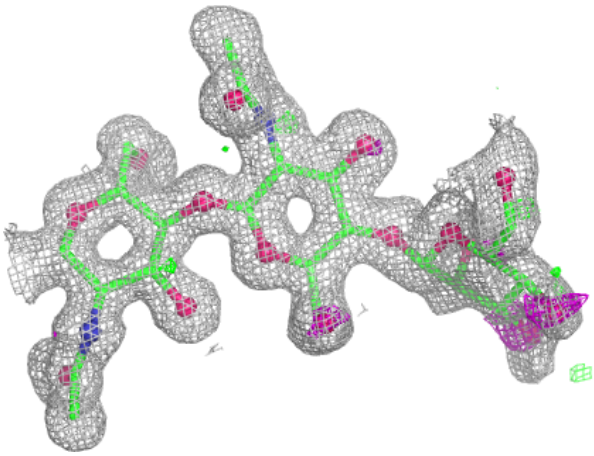


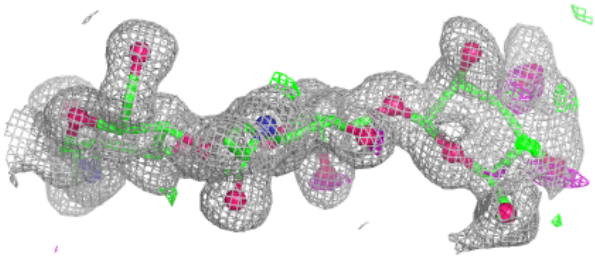


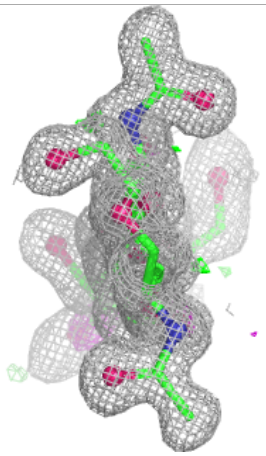


Electron density around Chain H:

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

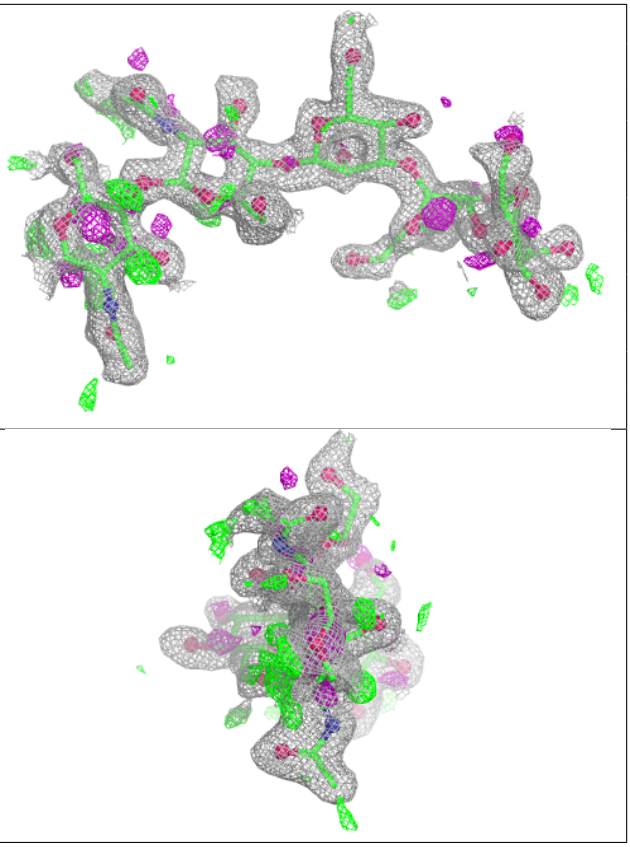






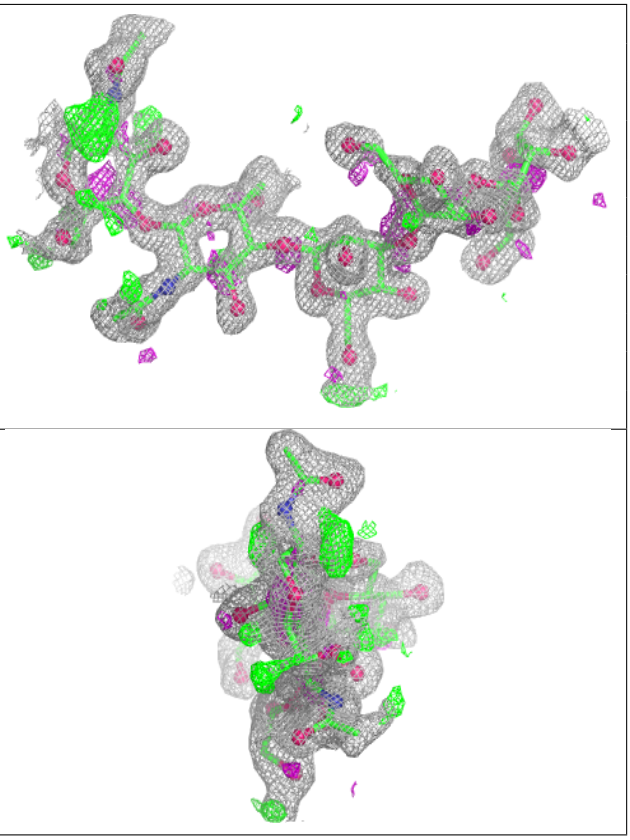
Electron density around Chain G:

$2mF_o$-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 L:

$2mF_o$-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	1005	14/15	0.71	0.29	27,30,31,35	0
8	SO4	A	1011	5/5	0.73	0.15	54,54,55,55	0
8	SO4	A	1012	5/5	0.81	0.14	53,53,54,54	0
10	GOL	A	1015[A]	6/6	0.81	0.22	22,24,24,25	6
10	GOL	A	1015[B]	6/6	0.81	0.22	23,24,25,27	6
8	SO4	B	1009	5/5	0.87	0.17	58,58,59,59	0
8	SO4	B	1008	5/5	0.88	0.12	43,43,44,45	0
10	GOL	B	1011[A]	6/6	0.88	0.17	24,24,24,24	6
10	GOL	B	1011[B]	6/6	0.88	0.17	23,24,24,26	6
5	NAG	A	1004	14/15	0.90	0.14	20,22,27,28	0
8	SO4	A	1013	5/5	0.90	0.25	54,55,55,55	0
5	NAG	B	1001	14/15	0.91	0.12	18,20,24,24	0
5	NAG	A	1001	14/15	0.92	0.10	16,19,22,22	0
5	NAG	B	1002	14/15	0.94	0.09	15,17,21,22	0
9	OXY	A	1014	2/2	0.95	0.24	24,24,24,25	0
9	OXY	B	1010	2/2	0.96	0.20	23,23,23,25	0
6	CU	A	1007	1/1	1.00	0.03	11,11,11,11	0
6	CU	A	1008	1/1	1.00	0.04	11,11,11,11	0
6	CU	A	1009	1/1	1.00	0.02	10,10,10,10	0
6	CU	B	1003	1/1	1.00	0.03	7,7,7,7	0
6	CU	B	1004	1/1	1.00	0.03	12,12,12,12	0
6	CU	B	1005	1/1	1.00	0.03	11,11,11,11	0
6	CU	B	1006	1/1	1.00	0.03	10,10,10,10	0
7	CL	A	1010	1/1	1.00	0.04	15,15,15,15	0
7	CL	B	1007	1/1	1.00	0.04	15,15,15,15	0
6	CU	A	1006	1/1	1.00	0.04	7,7,7,7	0

6.5 Other polymers (i)

There are no such residues in this entry.