



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 1, 2026 – 02:40 AM UTC

PDB ID : 4YMD / pdb_00004ymd
Title : CL-K1 trimer bound to man(alpha1-2)man
Authors : Wallis, R.; Venkatraman Girija, U.; Gingras, A.R.; Moody, P.C.E.; Marshall, J.E.
Deposited on : 2015-03-06
Resolution : 2.87 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 types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

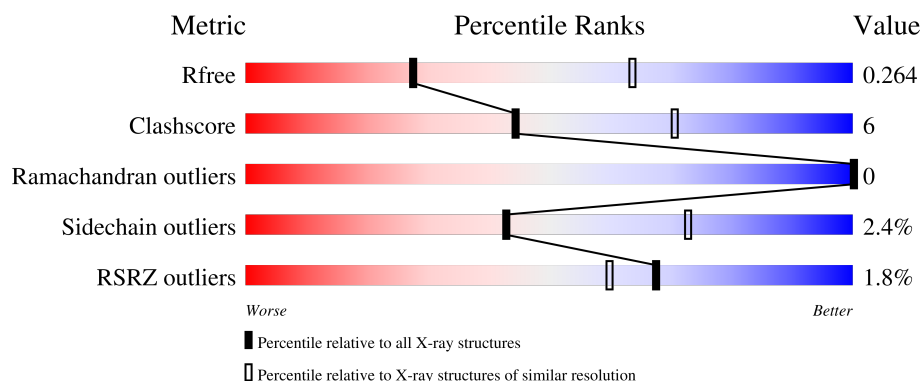
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

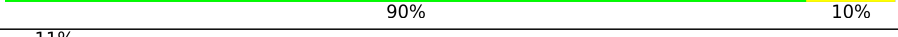
The reported resolution of this entry is 2.87 Å.

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}	180053	3557 (2.90-2.86)
Clashscore	190562	3801 (2.90-2.86)
Ramachandran outliers	187476	3699 (2.90-2.86)
Sidechain outliers	187428	3702 (2.90-2.86)
RSRZ outliers	180081	3558 (2.90-2.86)

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 $\leq 5\%$. 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	155	
1	B	155	
1	C	155	
1	D	155	
1	E	155	

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Mol	Chain	Length	Quality of chain
1	F	155	87% 11% ..
2	G	2	100%
2	H	2	50% 50%
2	I	2	50% 50%
2	J	2	50% 50%

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 7328 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 Collectin-11.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	154	Total	C	N	O	S	0	0	0
			1200	746	205	238	11			
1	B	154	Total	C	N	O	S	0	0	0
			1200	746	205	238	11			
1	D	155	Total	C	N	O	S	0	0	0
			1207	749	207	240	11			
1	F	154	Total	C	N	O	S	0	0	0
			1200	746	205	238	11			
1	C	153	Total	C	N	O	S	0	0	0
			1194	743	204	236	11			
1	E	149	Total	C	N	O	S	0	0	0
			1160	721	196	232	11			

- Molecule 2 is an oligosaccharide called alpha-D-mannopyranose-(1-2)-alpha-D-mannopyranose.

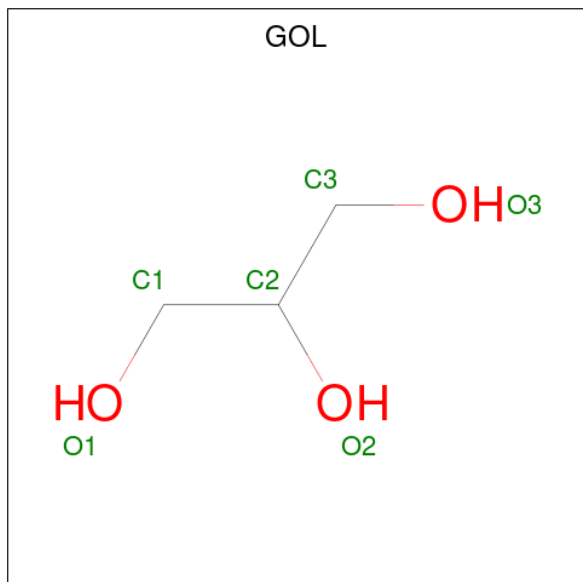


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace
2	G	2	Total	C	O	0	0	0
			23	12	11			
2	H	2	Total	C	O	0	0	0
			23	12	11			
2	I	2	Total	C	O	0	0	0
			23	12	11			
2	J	2	Total	C	O	0	0	0
			23	12	11			

- Molecule 3 is CALCIUM ION (CCD ID: CA) (formula: Ca).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	3	Total 3 Ca 3	0	0
3	B	3	Total 3 Ca 3	0	0
3	D	3	Total 3 Ca 3	0	0
3	F	3	Total 3 Ca 3	0	0
3	C	3	Total 3 Ca 3	0	0
3	E	3	Total 3 Ca 3	0	0

- Molecule 4 is GLYCEROL (CCD ID: GOL) (formula: $C_3H_8O_3$).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	1	Total 6 C 3 O 3	0	0
4	B	1	Total 6 C 3 O 3	0	0
4	D	1	Total 6 C 3 O 3	0	0
4	F	1	Total 6 C 3 O 3	0	0
4	C	1	Total 6 C 3 O 3	0	0
4	C	1	Total 6 C 3 O 3	0	0

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	E	1	Total	C	O	0	0
			6	3	3		

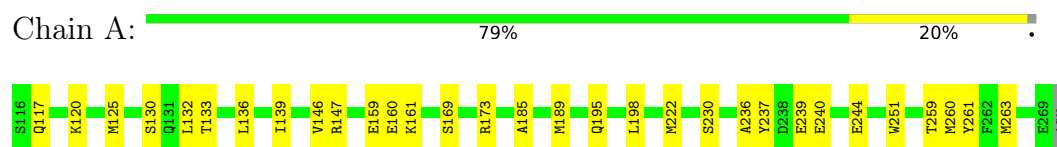
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	4	Total	O	0	0
			4	4		
5	B	3	Total	O	0	0
			3	3		
5	D	2	Total	O	0	0
			2	2		
5	F	3	Total	O	0	0
			3	3		
5	C	1	Total	O	0	0
			1	1		
5	E	2	Total	O	0	0
			2	2		

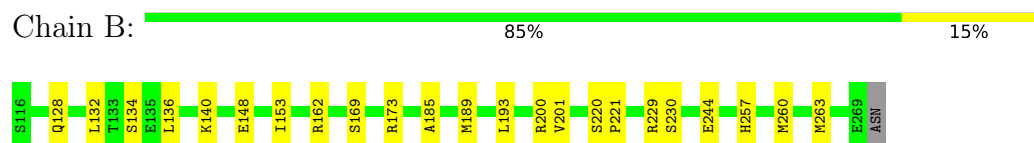
3 Residue-property plots [i](#)

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.

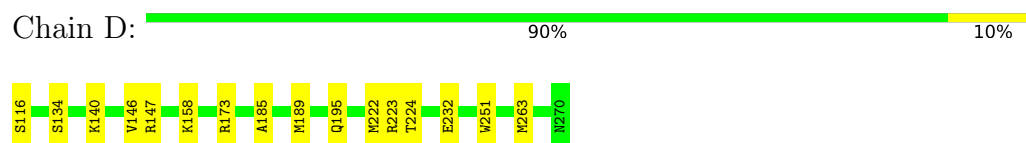
- Molecule 1: Collectin-11



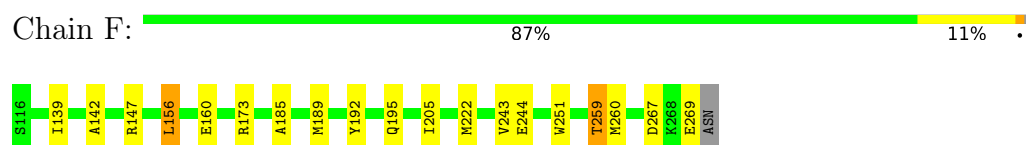
- Molecule 1: Collectin-11



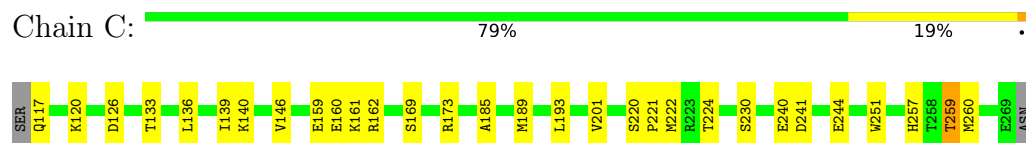
- Molecule 1: Collectin-11



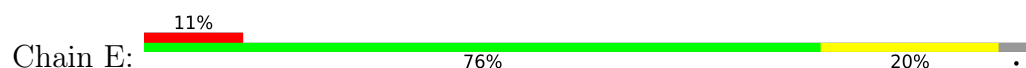
- Molecule 1: Collectin-11

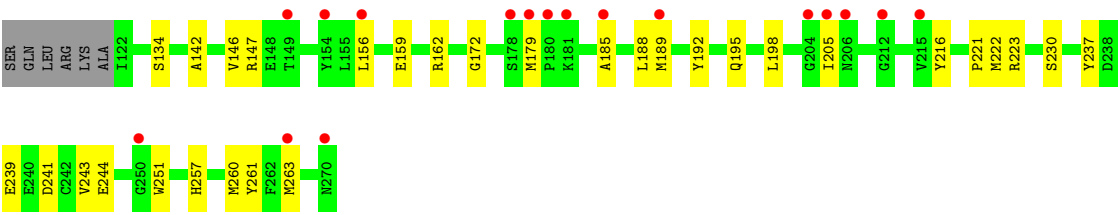


- Molecule 1: Collectin-11



- Molecule 1: Collectin-11





● Molecule 2: alpha-D-mannopyranose-(1-2)-alpha-D-mannopyranose

Chain G: 100%



● Molecule 2: alpha-D-mannopyranose-(1-2)-alpha-D-mannopyranose

Chain H: 50%



● Molecule 2: alpha-D-mannopyranose-(1-2)-alpha-D-mannopyranose

Chain I: 50%



● Molecule 2: alpha-D-mannopyranose-(1-2)-alpha-D-mannopyranose

Chain J: 50%



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	71.85Å 106.65Å 78.39Å 90.00° 92.17° 90.00°	Depositor
Resolution (Å)	44.08 – 2.87 44.08 – 2.87	Depositor EDS
% Data completeness (in resolution range)	91.0 (44.08-2.87) 91.0 (44.08-2.87)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.55 (at 2.86Å)	Xtriage
Refinement program	PHENIX	Depositor
R, R_{free}	0.213 , 0.264 0.217 , 0.264	Depositor DCC
R_{free} test set	1232 reflections (4.54%)	wwPDB-VP
Wilson B-factor (Å ²)	79.8	Xtriage
Anisotropy	0.136	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 61.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	0.037 for h,-k,-l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	7328	wwPDB-VP
Average B, all atoms (Å ²)	95.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.57% 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: CA, MAN, GOL

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	0.31	0/1220	0.75	0/1640
1	B	0.30	0/1220	0.74	0/1640
1	C	0.37	0/1214	0.78	0/1632
1	D	0.30	0/1225	0.75	0/1647
1	E	0.31	0/1180	0.77	0/1588
1	F	0.31	0/1220	0.75	0/1640
All	All	0.32	0/7279	0.76	0/9787

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [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	1200	0	1149	18	0
1	B	1200	0	1149	13	0
1	C	1194	0	1144	18	0
1	D	1207	0	1150	16	0
1	E	1160	0	1100	18	0
1	F	1200	0	1149	12	0
2	G	23	0	19	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	H	23	0	19	2	0
2	I	23	0	19	1	0
2	J	23	0	19	0	0
3	A	3	0	0	0	0
3	B	3	0	0	0	0
3	C	3	0	0	0	0
3	D	3	0	0	0	0
3	E	3	0	0	0	0
3	F	3	0	0	0	0
4	A	6	0	7	0	0
4	B	6	0	6	0	0
4	C	12	0	13	2	0
4	D	6	0	6	0	0
4	E	6	0	8	1	0
4	F	6	0	7	0	0
5	A	4	0	0	0	0
5	B	3	0	0	0	0
5	C	1	0	0	0	0
5	D	2	0	0	1	0
5	E	2	0	0	0	0
5	F	3	0	0	0	0
All	All	7328	0	6964	85	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 6.

All (85) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:223:ARG:NH2	1:D:223:ARG:NH1	2.29	0.80
1:D:223:ARG:NH1	1:D:223:ARG:NE	2.30	0.79
1:D:223:ARG:NH2	1:D:223:ARG:NE	2.30	0.79
1:D:173:ARG:NH1	1:F:142:ALA:O	2.18	0.77
1:B:229:ARG:HH22	2:H:2:MAN:H3	1.54	0.72
1:B:169:SER:OG	1:B:173:ARG:NH1	2.26	0.69
1:B:200:ARG:NH2	2:H:2:MAN:O3	2.26	0.67
1:A:147:ARG:NH2	1:A:195:GLN:OE1	2.30	0.64
1:F:222:MET:HE1	1:F:251:TRP:HZ3	1.64	0.63
1:C:169:SER:OG	1:C:173:ARG:NH1	2.33	0.61
1:A:133:THR:HG22	1:B:132:LEU:HD11	1.82	0.61
1:D:222:MET:HE1	1:D:251:TRP:HZ3	1.63	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:147:ARG:NH2	1:F:195:GLN:OE1	2.23	0.61
1:C:159:GLU:OE2	1:C:161:LYS:NZ	2.33	0.60
1:E:185:ALA:O	1:E:189:MET:HG3	2.02	0.59
1:C:222:MET:HE1	1:C:251:TRP:HZ3	1.68	0.59
1:A:169:SER:OG	1:A:173:ARG:NH2	2.36	0.59
1:F:173:ARG:NH1	1:E:142:ALA:O	2.37	0.58
1:C:117:GLN:HA	1:C:120:LYS:HE3	1.86	0.58
1:A:222:MET:HE1	1:A:251:TRP:HZ3	1.69	0.58
1:E:222:MET:HE1	1:E:251:TRP:HZ3	1.70	0.57
1:B:162:ARG:HD2	1:B:257:HIS:HA	1.87	0.56
1:D:158:LYS:NZ	1:E:159:GLU:OE2	2.31	0.56
1:A:185:ALA:O	1:A:189:MET:HG3	2.07	0.55
1:B:185:ALA:O	1:B:189:MET:HG3	2.06	0.55
1:F:267:ASP:HB3	1:F:269:GLU:OE1	2.07	0.54
1:C:240:GLU:OE2	4:C:301:GOL:O2	2.27	0.53
1:C:162:ARG:HD2	1:C:257:HIS:HA	1.90	0.52
1:A:237:TYR:HB2	1:A:239:GLU:HG2	1.92	0.52
1:F:185:ALA:O	1:F:189:MET:HG3	2.10	0.51
1:D:222:MET:HE1	1:D:251:TRP:CZ3	2.44	0.50
1:F:222:MET:HE1	1:F:251:TRP:CZ3	2.46	0.50
1:C:185:ALA:O	1:C:189:MET:HG3	2.11	0.50
1:A:130:SER:O	1:A:133:THR:OG1	2.25	0.50
1:A:222:MET:HE1	1:A:251:TRP:CZ3	2.46	0.50
1:E:156:LEU:HD12	1:E:263:MET:HG3	1.93	0.50
1:E:237:TYR:HB2	1:E:239:GLU:HG2	1.92	0.50
1:D:185:ALA:O	1:D:189:MET:HG3	2.11	0.50
1:E:241:ASP:OD2	4:E:304:GOL:O2	2.23	0.49
1:D:189:MET:HG2	1:D:263:MET:HE1	1.93	0.48
1:C:160:GLU:HB3	1:C:259:THR:CG2	2.44	0.48
1:A:189:MET:HG2	1:A:263:MET:HE1	1.94	0.48
1:C:241:ASP:OD2	4:C:305:GOL:O1	2.32	0.48
1:B:140:LYS:HG2	1:C:139:ILE:HD11	1.96	0.47
1:A:160:GLU:HB3	1:A:259:THR:CG2	2.44	0.47
1:B:193:LEU:HD11	1:B:201:VAL:HG22	1.97	0.47
1:E:222:MET:HE1	1:E:251:TRP:CZ3	2.49	0.46
1:D:195:GLN:NE2	1:E:172:GLY:HA3	2.30	0.46
1:E:156:LEU:HD13	1:E:192:TYR:CE2	2.51	0.46
1:E:189:MET:HG2	1:E:263:MET:HE1	1.98	0.46
1:D:232:GLU:OE2	2:I:1:MAN:O4	2.34	0.46
1:F:160:GLU:HB3	1:F:259:THR:CG2	2.46	0.46
1:F:244:GLU:HB3	1:F:260:MET:HE1	1.97	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:189:MET:HG2	1:B:263:MET:HE1	1.99	0.45
1:B:244:GLU:HB3	1:B:260:MET:HE1	1.99	0.45
1:B:220:SER:HA	1:B:221:PRO:HD3	1.87	0.44
1:A:139:ILE:HD11	1:C:140:LYS:HG2	1.99	0.44
1:C:244:GLU:HB3	1:C:260:MET:HE1	2.00	0.43
1:E:198:LEU:HD21	1:E:261:TYR:HD2	1.83	0.43
1:C:140:LYS:O	1:C:146:VAL:HG22	2.18	0.43
1:C:193:LEU:HD11	1:C:201:VAL:HG22	2.01	0.43
1:B:148:GLU:HG3	1:B:153:ILE:HG12	2.01	0.42
1:D:116:SER:N	5:D:401:HOH:O	2.52	0.42
1:B:128:GLN:O	1:B:132:LEU:HD13	2.19	0.42
1:A:159:GLU:OE1	1:A:161:LYS:NZ	2.53	0.42
1:A:117:GLN:HA	1:A:120:LYS:HE3	2.02	0.42
1:D:223:ARG:NH1	1:D:223:ARG:CD	2.82	0.42
1:C:220:SER:HA	1:C:221:PRO:HD3	1.83	0.42
1:E:147:ARG:NH2	1:E:195:GLN:OE1	2.50	0.42
1:F:156:LEU:HD22	1:F:192:TYR:CE2	2.55	0.41
1:D:140:LYS:O	1:D:146:VAL:HG22	2.20	0.41
1:A:125:MET:HE1	1:C:126:ASP:HA	2.02	0.41
1:A:132:LEU:HD11	1:C:133:THR:HG23	2.02	0.41
1:D:147:ARG:NH2	1:D:195:GLN:OE1	2.53	0.41
1:A:236:ALA:HB3	1:A:240:GLU:CD	2.46	0.41
1:C:222:MET:HE1	1:C:251:TRP:CZ3	2.52	0.41
1:E:179:MET:HB3	1:E:216:TYR:HD2	1.85	0.41
1:E:221:PRO:O	1:E:223:ARG:NH1	2.54	0.41
1:D:140:LYS:HG2	1:F:139:ILE:HD11	2.03	0.40
1:E:244:GLU:HB3	1:E:260:MET:HE1	2.03	0.40
1:A:244:GLU:HB3	1:A:260:MET:HE1	2.02	0.40
1:F:205:ILE:HB	1:F:243:VAL:HB	2.03	0.40
1:E:205:ILE:HB	1:E:243:VAL:HB	2.04	0.40
1:A:198:LEU:HD21	1:A:261:TYR:HD2	1.87	0.40
1:E:162:ARG:HD2	1:E:257:HIS:HA	2.03	0.40

There are no symmetry-related clashes.

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	152/155 (98%)	146 (96%)	6 (4%)	0	100	100
1	B	152/155 (98%)	145 (95%)	7 (5%)	0	100	100
1	C	151/155 (97%)	145 (96%)	6 (4%)	0	100	100
1	D	153/155 (99%)	147 (96%)	6 (4%)	0	100	100
1	E	147/155 (95%)	142 (97%)	5 (3%)	0	100	100
1	F	152/155 (98%)	146 (96%)	6 (4%)	0	100	100
All	All	907/930 (98%)	871 (96%)	36 (4%)	0	100	100

There are no Ramachandran outliers to report.

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	126/127 (99%)	123 (98%)	3 (2%)	43	73
1	B	126/127 (99%)	123 (98%)	3 (2%)	43	73
1	C	125/127 (98%)	121 (97%)	4 (3%)	34	65
1	D	126/127 (99%)	124 (98%)	2 (2%)	55	81
1	E	122/127 (96%)	118 (97%)	4 (3%)	33	65
1	F	126/127 (99%)	124 (98%)	2 (2%)	55	81
All	All	751/762 (99%)	733 (98%)	18 (2%)	43	73

All (18) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	136	LEU
1	A	146	VAL
1	A	230	SER

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Mol	Chain	Res	Type
1	B	134	SER
1	B	136	LEU
1	B	230	SER
1	D	134	SER
1	D	224	THR
1	F	156	LEU
1	F	259	THR
1	C	136	LEU
1	C	224	THR
1	C	230	SER
1	C	259	THR
1	E	134	SER
1	E	146	VAL
1	E	188	LEU
1	E	230	SER

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

Mol	Chain	Res	Type
1	B	235	ASN
1	F	219	HIS
1	F	235	ASN
1	C	235	ASN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

8 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	MAN	G	1	2,3	12,12,12	0.51	0	17,17,17	1.00	1 (5%)
2	MAN	G	2	2	11,11,12	1.03	1 (9%)	15,15,17	1.76	3 (20%)
2	MAN	H	1	2,3	12,12,12	0.60	0	17,17,17	1.02	1 (5%)
2	MAN	H	2	2	11,11,12	1.05	2 (18%)	15,15,17	1.16	0
2	MAN	I	1	2,3	12,12,12	0.46	0	17,17,17	1.13	1 (5%)
2	MAN	I	2	2	11,11,12	0.74	0	15,15,17	0.88	1 (6%)
2	MAN	J	1	2,3	12,12,12	0.47	0	17,17,17	1.08	1 (5%)
2	MAN	J	2	2	11,11,12	0.84	0	15,15,17	0.88	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
2	MAN	G	1	2,3	-	2/2/22/22	0/1/1/1
2	MAN	G	2	2	-	0/2/19/22	0/1/1/1
2	MAN	H	1	2,3	-	0/2/22/22	0/1/1/1
2	MAN	H	2	2	-	0/2/19/22	0/1/1/1
2	MAN	I	1	2,3	-	2/2/22/22	0/1/1/1
2	MAN	I	2	2	-	0/2/19/22	0/1/1/1
2	MAN	J	1	2,3	-	0/2/22/22	0/1/1/1
2	MAN	J	2	2	-	1/2/19/22	0/1/1/1

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	G	2	MAN	C1-C2	2.84	1.59	1.52
2	H	2	MAN	C2-C3	2.32	1.56	1.52
2	H	2	MAN	C1-C2	2.12	1.57	1.52

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	G	2	MAN	C1-O5-C5	3.73	117.18	112.19
2	G	2	MAN	C1-C2-C3	3.39	114.58	109.64

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	G	1	MAN	O2-C2-C3	-3.07	103.14	110.38
2	H	1	MAN	O2-C2-C3	-3.03	103.24	110.38
2	I	1	MAN	O2-C2-C3	-2.95	103.42	110.38
2	G	2	MAN	O5-C1-C2	2.87	117.64	110.79
2	J	1	MAN	O2-C2-C3	-2.82	103.73	110.38
2	I	2	MAN	O2-C2-C3	-2.04	105.93	110.15

There are no chirality outliers.

All (5) torsion outliers are listed below:

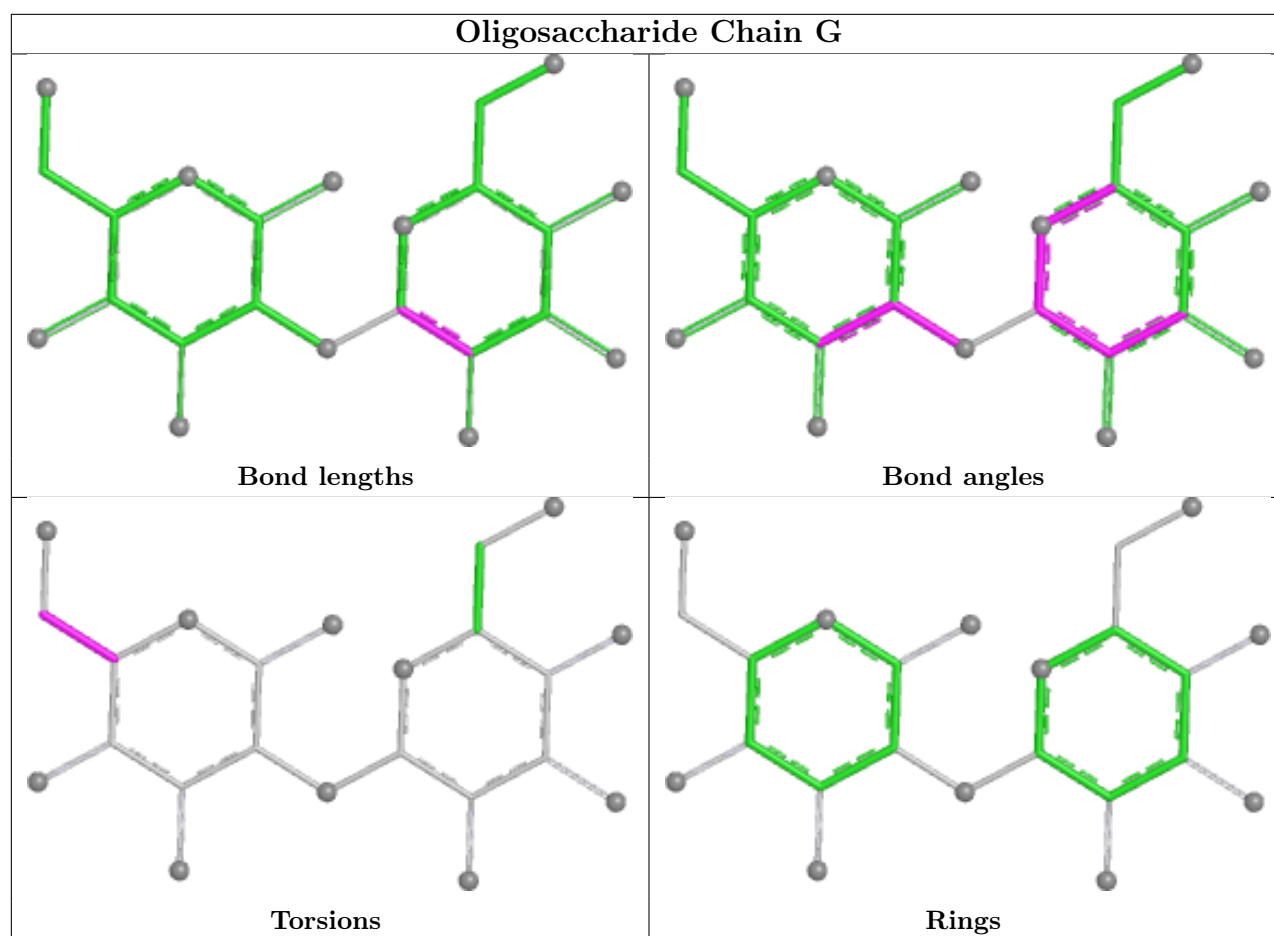
Mol	Chain	Res	Type	Atoms
2	G	1	MAN	O5-C5-C6-O6
2	G	1	MAN	C4-C5-C6-O6
2	I	1	MAN	C4-C5-C6-O6
2	I	1	MAN	O5-C5-C6-O6
2	J	2	MAN	C4-C5-C6-O6

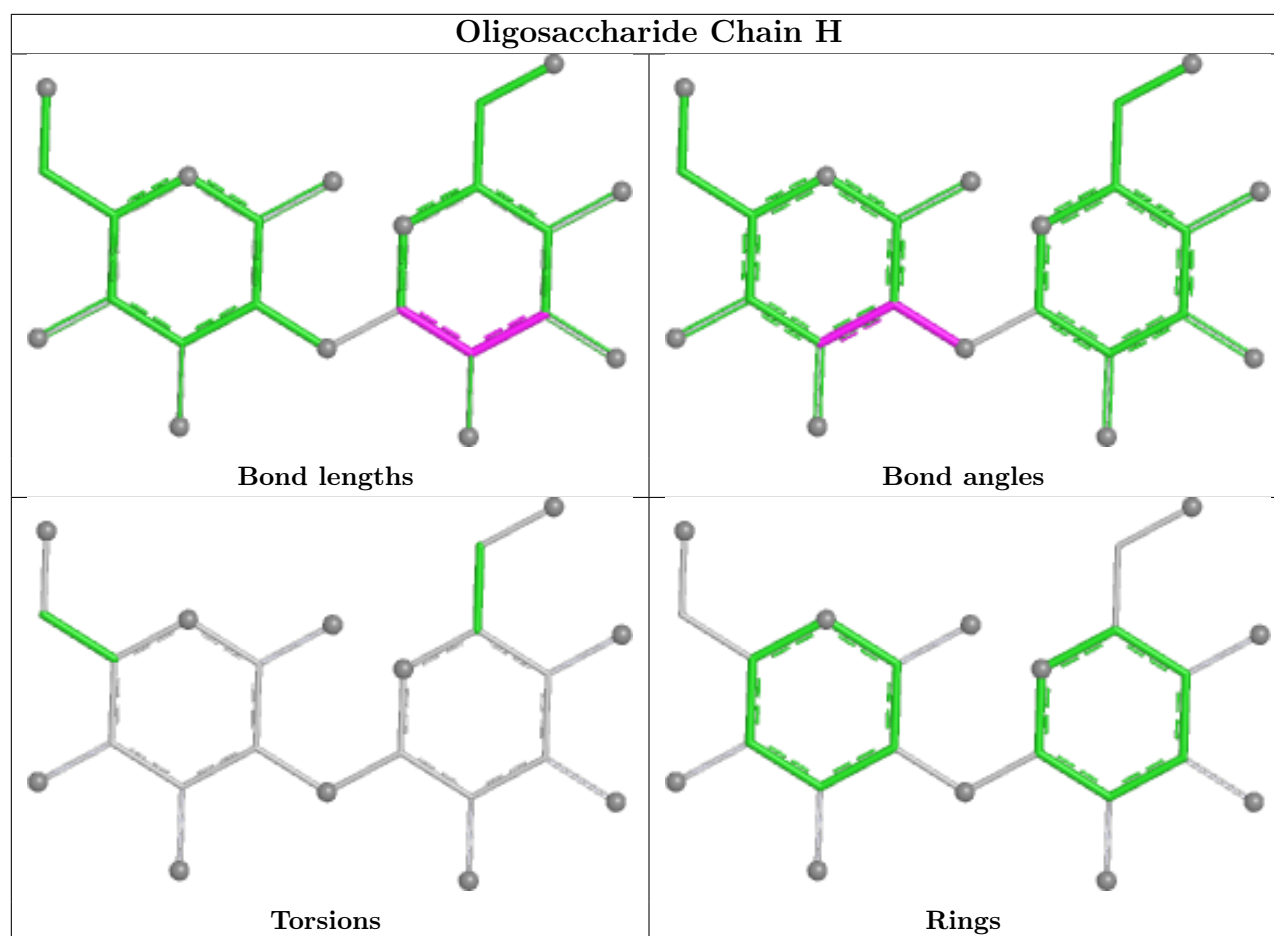
There are no ring outliers.

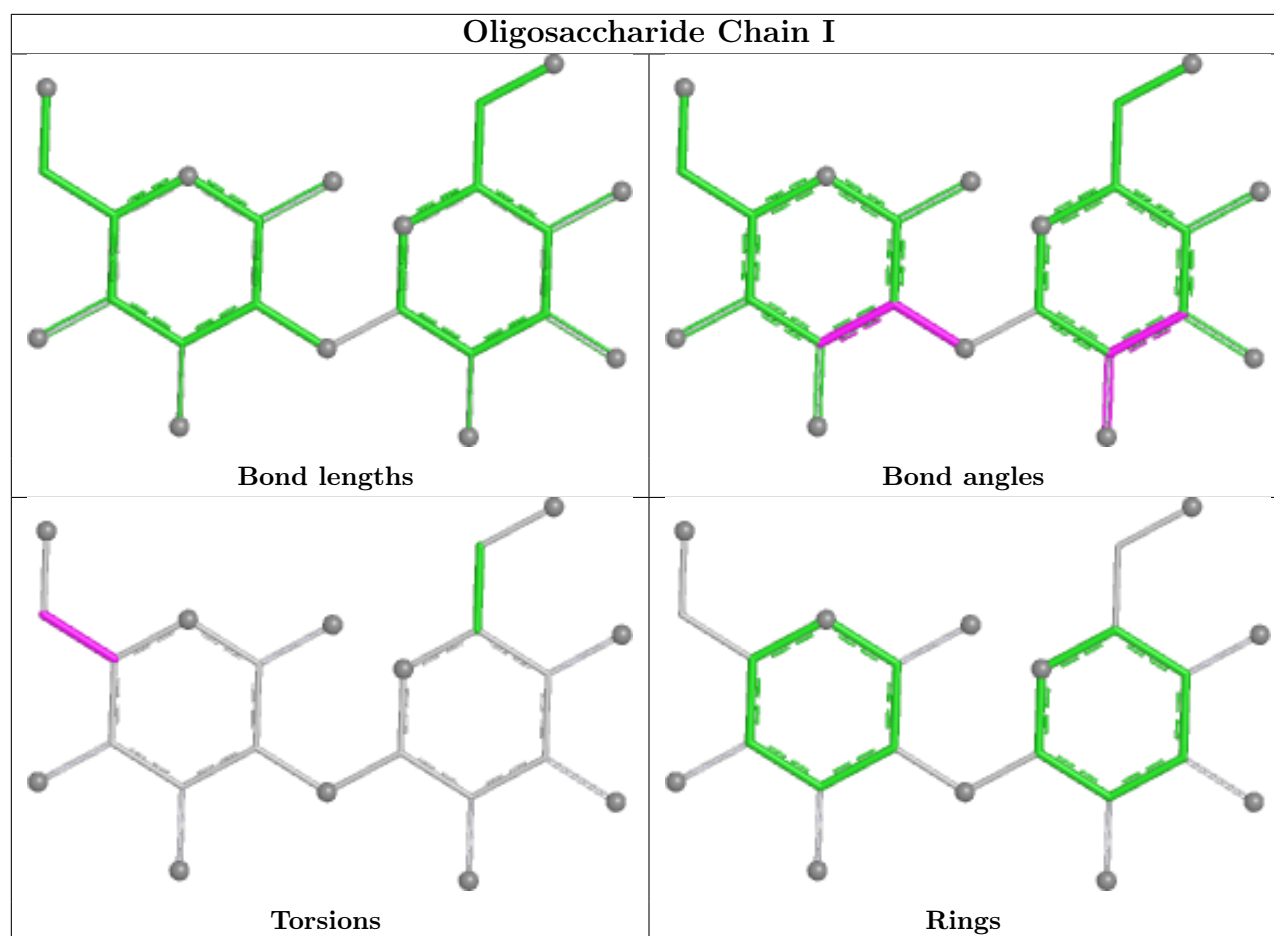
2 monomers are involved in 3 short contacts:

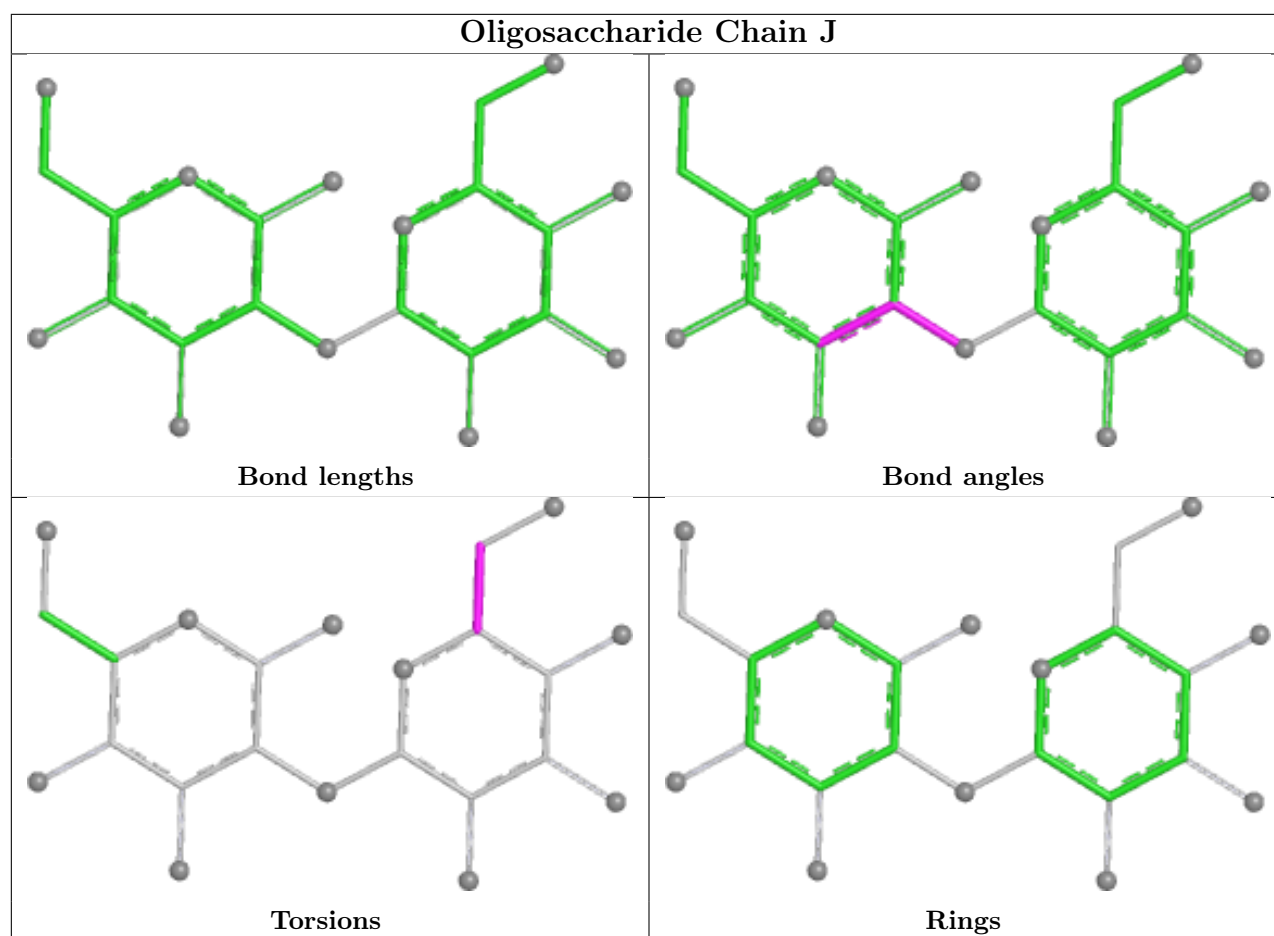
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	I	1	MAN	1	0
2	H	2	MAN	2	0

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









5.6 Ligand geometry [i](#)

Of 25 ligands modelled in this entry, 18 are monoatomic - leaving 7 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	$\# Z > 2$
4	GOL	C	301	3	5,5,5	0.42	0	5,5,5	0.31	0
4	GOL	F	306	3	5,5,5	0.37	0	5,5,5	0.31	0
4	GOL	A	306	3	5,5,5	0.37	0	5,5,5	0.26	0
4	GOL	C	305	3	5,5,5	0.38	0	5,5,5	0.31	0
4	GOL	B	306	3	5,5,5	0.38	0	5,5,5	0.33	0
4	GOL	D	306	3	5,5,5	0.36	0	5,5,5	0.38	0
4	GOL	E	304	3	5,5,5	0.39	0	5,5,5	0.21	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
4	GOL	C	301	3	-	1/4/4/4	-
4	GOL	F	306	3	-	2/4/4/4	-
4	GOL	A	306	3	-	2/4/4/4	-
4	GOL	C	305	3	-	4/4/4/4	-
4	GOL	B	306	3	-	4/4/4/4	-
4	GOL	D	306	3	-	2/4/4/4	-
4	GOL	E	304	3	-	0/4/4/4	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (15) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	306	GOL	O1-C1-C2-C3
4	B	306	GOL	O1-C1-C2-C3
4	B	306	GOL	C1-C2-C3-O3
4	D	306	GOL	O1-C1-C2-C3
4	F	306	GOL	O1-C1-C2-C3
4	C	305	GOL	C1-C2-C3-O3
4	A	306	GOL	O1-C1-C2-O2
4	D	306	GOL	O1-C1-C2-O2
4	B	306	GOL	O1-C1-C2-O2
4	F	306	GOL	O1-C1-C2-O2
4	B	306	GOL	O2-C2-C3-O3
4	C	305	GOL	O2-C2-C3-O3
4	C	305	GOL	O1-C1-C2-O2
4	C	305	GOL	O1-C1-C2-C3
4	C	301	GOL	O1-C1-C2-O2

There are no ring outliers.

3 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	C	301	GOL	1	0
4	C	305	GOL	1	0

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Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	E	304	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

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	154/155 (99%)	-0.20	0	100	100	56, 88, 143, 233	0
1	B	154/155 (99%)	-0.40	0	100	100	50, 76, 106, 148	0
1	C	153/155 (98%)	-0.25	0	100	100	63, 98, 136, 159	0
1	D	155/155 (100%)	-0.41	0	100	100	52, 80, 118, 157	0
1	E	149/155 (96%)	0.67	17 (11%)	10	8	64, 139, 199, 245	0
1	F	154/155 (99%)	-0.45	0	100	100	50, 73, 103, 121	0
All	All	919/930 (98%)	-0.18	17 (1%)	67	60	50, 85, 159, 245	0

All (17) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	E	205	ILE	4.1
1	E	156	LEU	3.0
1	E	185	ALA	2.8
1	E	212	GLY	2.7
1	E	204	GLY	2.6
1	E	178	SER	2.5
1	E	270	ASN	2.5
1	E	189	MET	2.5
1	E	250	GLY	2.5
1	E	154	TYR	2.5
1	E	206	ASN	2.4
1	E	215	VAL	2.4
1	E	180	PRO	2.1
1	E	149	THR	2.1
1	E	181	LYS	2.1
1	E	179	MET	2.1
1	E	263	MET	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 ⓘ

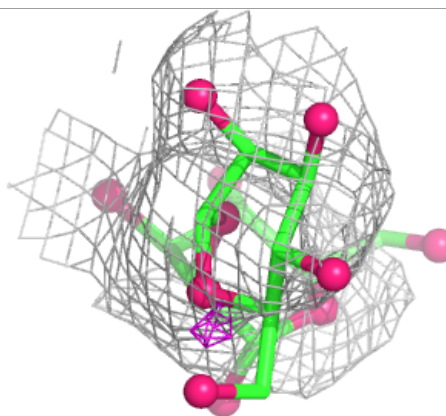
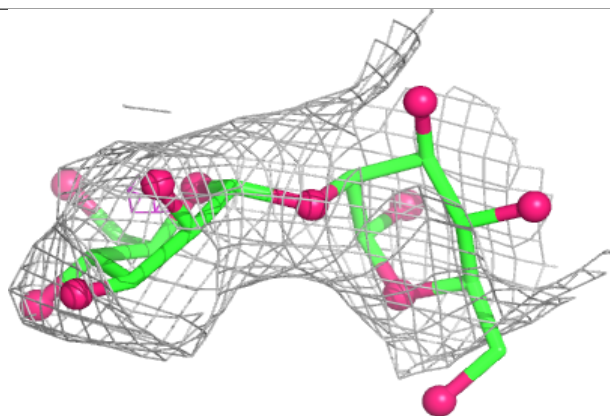
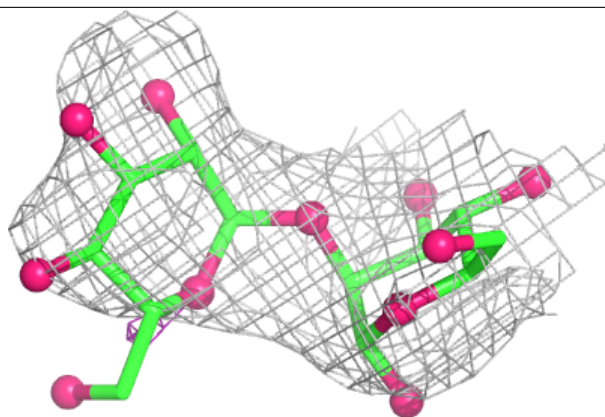
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	MAN	H	2	11/12	0.68	0.13	116,123,132,136	0
2	MAN	I	2	11/12	0.69	0.10	127,136,140,140	0
2	MAN	J	2	11/12	0.71	0.12	132,136,148,152	0
2	MAN	G	2	11/12	0.76	0.10	93,123,128,132	0
2	MAN	H	1	12/12	0.86	0.12	57,102,118,121	0
2	MAN	I	1	12/12	0.87	0.09	81,103,124,130	0
2	MAN	J	1	12/12	0.88	0.10	94,117,126,128	0
2	MAN	G	1	12/12	0.93	0.08	84,104,119,121	0

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

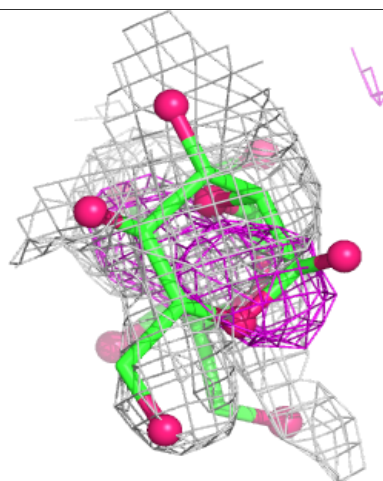
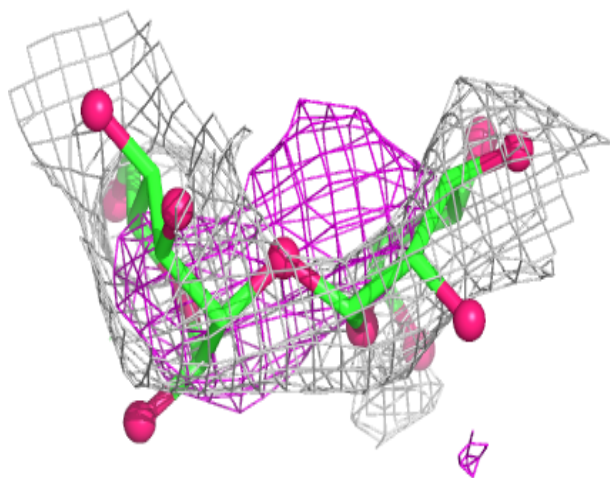
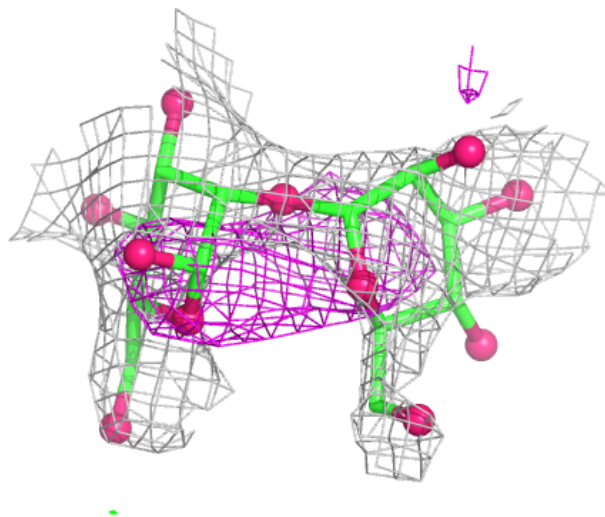
Electron density around Chain G:

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



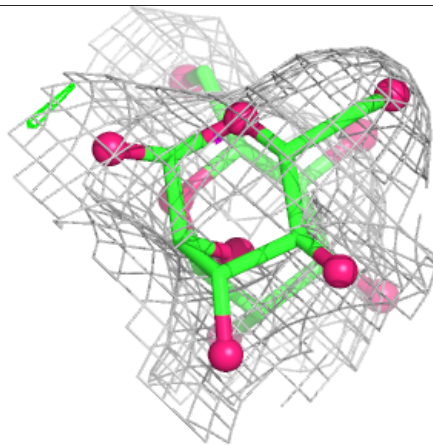
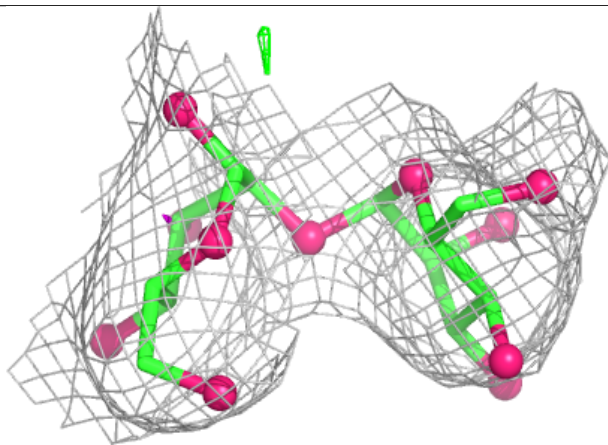
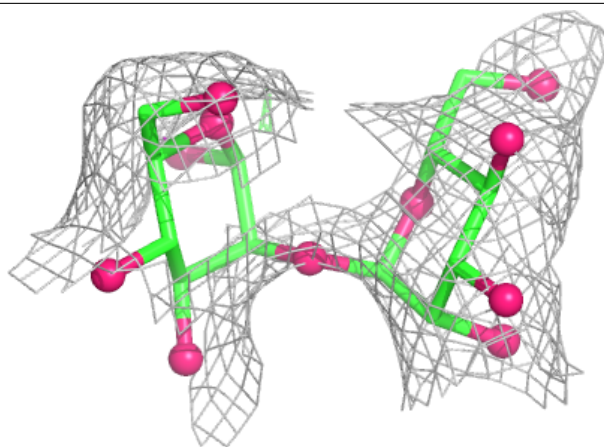
Electron density around Chain H:

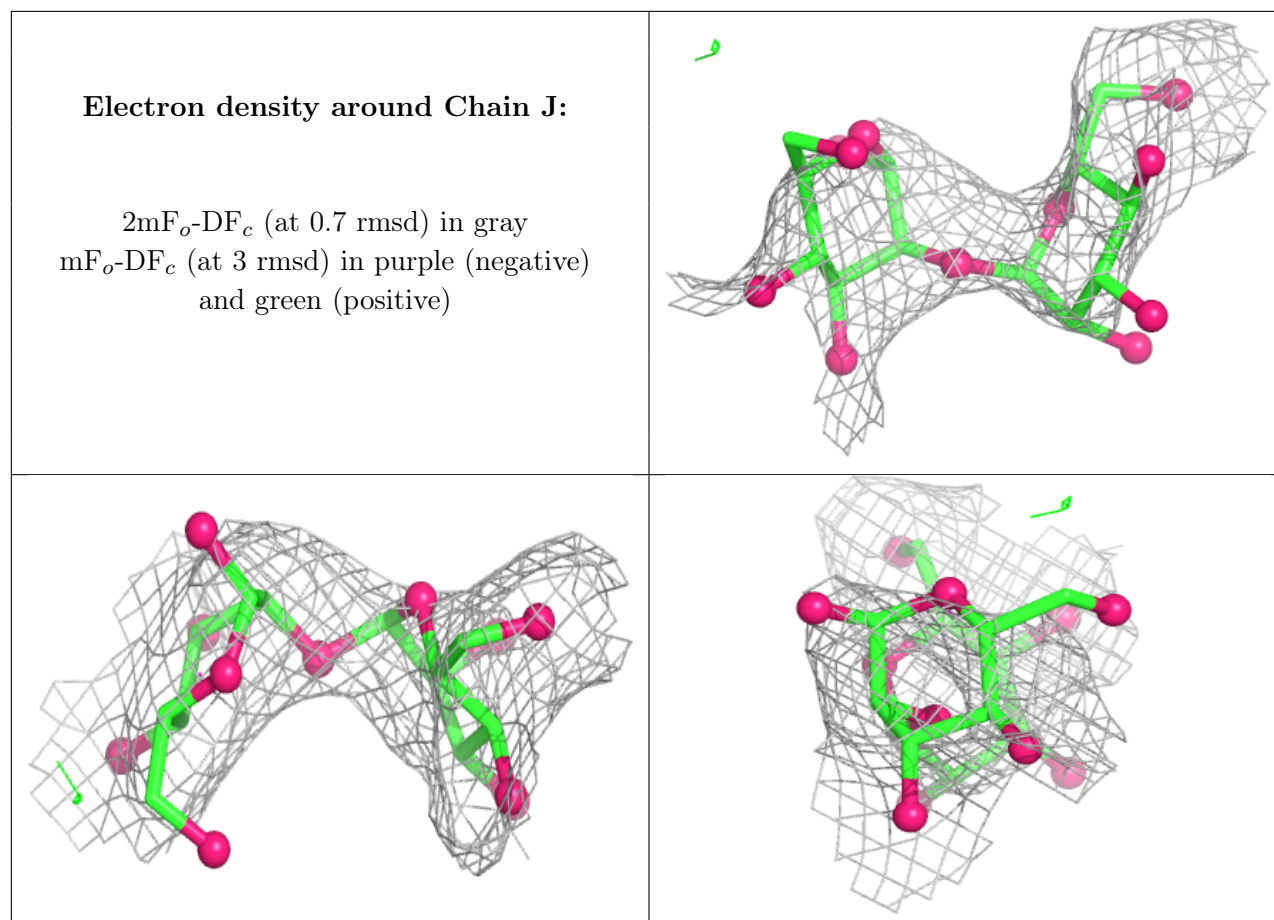
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around Chain I:

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





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
3	CA	E	301	1/1	0.69	0.08	173,173,173,173	0
4	GOL	A	306	6/6	0.80	0.15	88,102,107,108	0
4	GOL	D	306	6/6	0.83	0.13	47,85,89,98	0
3	CA	E	303	1/1	0.84	0.07	184,184,184,184	0
4	GOL	F	306	6/6	0.85	0.11	60,71,77,78	0
4	GOL	C	305	6/6	0.87	0.07	94,106,114,119	0
4	GOL	B	306	6/6	0.91	0.15	43,70,75,78	0
4	GOL	E	304	6/6	0.91	0.07	96,100,102,102	0
4	GOL	C	301	6/6	0.93	0.09	55,72,75,82	0
3	CA	C	304	1/1	0.95	0.05	139,139,139,139	0
3	CA	F	305	1/1	0.96	0.07	105,105,105,105	0
3	CA	E	302	1/1	0.97	0.03	140,140,140,140	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	CA	F	303	1/1	0.97	0.09	72,72,72,72	0
3	CA	A	305	1/1	0.97	0.07	102,102,102,102	0
3	CA	C	303	1/1	0.98	0.03	87,87,87,87	0
3	CA	A	304	1/1	0.98	0.03	73,73,73,73	0
3	CA	B	305	1/1	0.98	0.07	82,82,82,82	0
3	CA	D	304	1/1	0.99	0.04	72,72,72,72	0
3	CA	D	305	1/1	0.99	0.03	71,71,71,71	0
3	CA	B	304	1/1	0.99	0.02	60,60,60,60	0
3	CA	F	304	1/1	0.99	0.05	71,71,71,71	0
3	CA	D	303	1/1	0.99	0.02	58,58,58,58	0
3	CA	C	302	1/1	0.99	0.03	105,105,105,105	0
3	CA	B	303	1/1	1.00	0.04	59,59,59,59	0
3	CA	A	303	1/1	1.00	0.06	78,78,78,78	0

6.5 Other polymers [i](#)

There are no such residues in this entry.