



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 03:03 PM UTC

PDB ID : 7BTL / pdb_00007btl
Title : Mevo lectin complex with mannopentose
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Deposited on : 2020-04-01
Resolution : 2.25 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
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
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

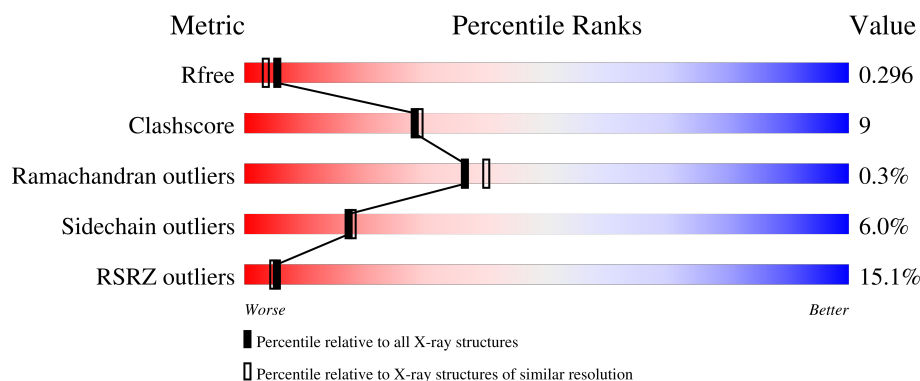
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.25 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	1898 (2.26-2.26)
Clashscore	190562	2005 (2.26-2.26)
Ramachandran outliers	187476	1965 (2.26-2.26)
Sidechain outliers	187428	1966 (2.26-2.26)
RSRZ outliers	180081	1898 (2.26-2.26)

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	140	2% 81% 17% .
1	B	140	4% 80% 18% .
1	C	140	17% 79% 17% .
1	D	140	9% 78% 20% .
1	E	140	30% 78% 21% .

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Mol	Chain	Length	Quality of chain
1	F	140	
1	G	140	
2	H	4	
2	I	4	
2	M	4	
3	J	3	
4	K	2	
4	L	2	

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	MAN	G	201	-	-	X	-

2 Entry composition [i](#)

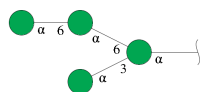
There are 7 unique types of molecules in this entry. The entry contains 7802 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 lectin.

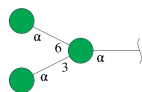
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	B	140	Total	C	N	O	S	0	0	0
			1065	678	173	211	3			
1	G	139	Total	C	N	O	S	0	0	0
			1034	660	166	205	3			
1	A	140	Total	C	N	O	S	0	0	0
			1059	676	172	208	3			
1	D	140	Total	C	N	O	S	0	0	0
			1058	675	172	208	3			
1	E	140	Total	C	N	O	S	0	0	0
			1065	679	174	209	3			
1	F	140	Total	C	N	O	S	0	0	0
			1061	676	173	209	3			
1	C	140	Total	C	N	O	S	0	0	0
			1053	672	170	208	3			

- Molecule 2 is an oligosaccharide called alpha-D-mannopyranose-(1-6)-alpha-D-mannopyranose-(1-6)-[alpha-D-mannopyranose-(1-3)]alpha-D-mannopyranose.



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace
2	H	4	Total	C	O	0	0	0
			45	24	21			
2	I	4	Total	C	O	0	0	0
			45	24	21			
2	M	4	Total	C	O	0	0	0
			45	24	21			

- Molecule 3 is an oligosaccharide called alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-6)]alpha-D-mannopyranose.



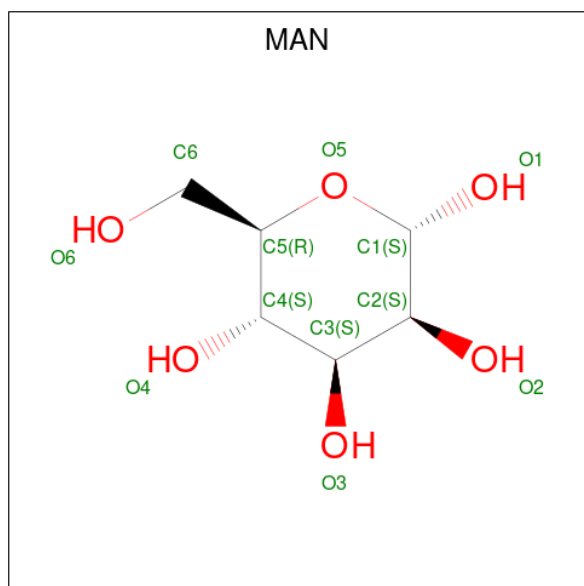
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace
3	J	3	Total	C	O	0	0	0
			34	18	16			

- Molecule 4 is an oligosaccharide called alpha-D-mannopyranose-(1-6)-alpha-D-mannopyranose.



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace
4	K	2	Total	C	O	0	0	0
			23	12	11			
4	L	2	Total	C	O	0	0	0
			23	12	11			

- Molecule 5 is alpha-D-mannopyranose (CCD ID: MAN) (formula: C₆H₁₂O₆) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	G	1	Total	C	O	0	0
			12	6	6		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	D	1	Total	C	O	0	0
			11	6	5		

- Molecule 6 is GLYCEROL (CCD ID: GOL) (formula: $C_3H_8O_3$).



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

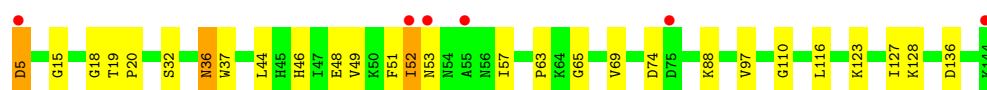
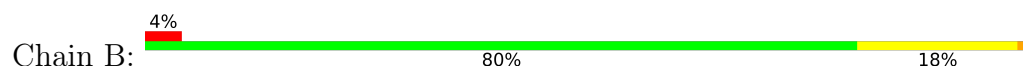
- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	B	33	Total	O	0	0
			33	33		
7	G	18	Total	O	0	0
			18	18		
7	A	27	Total	O	0	0
			27	27		
7	D	15	Total	O	0	0
			15	15		
7	E	31	Total	O	0	0
			31	31		
7	F	22	Total	O	0	0
			22	22		
7	C	17	Total	O	0	0
			17	17		

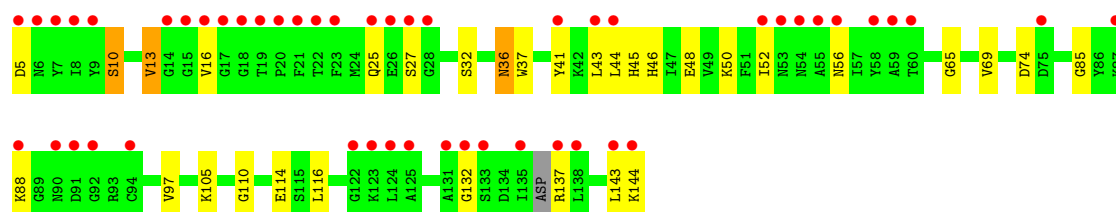
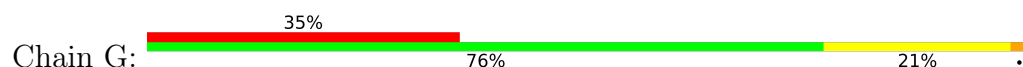
3 Residue-property plots [i](#)

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.

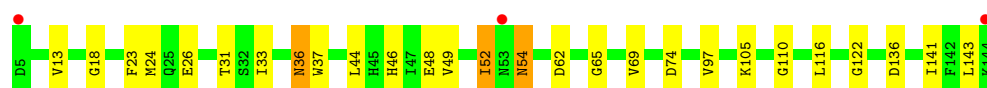
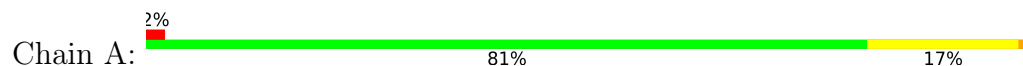
- Molecule 1: lectin



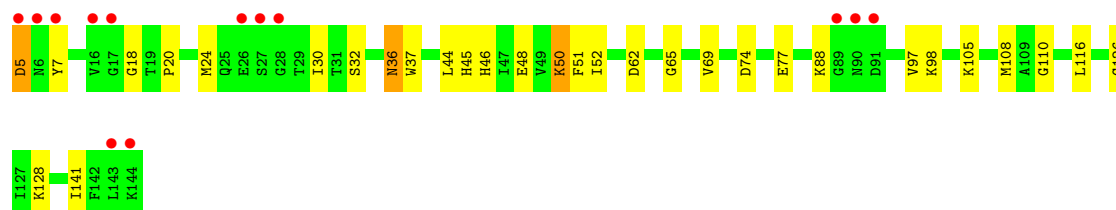
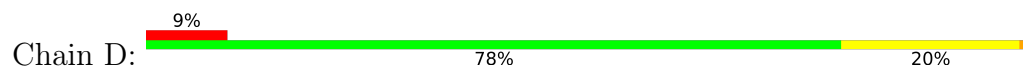
- Molecule 1: lectin



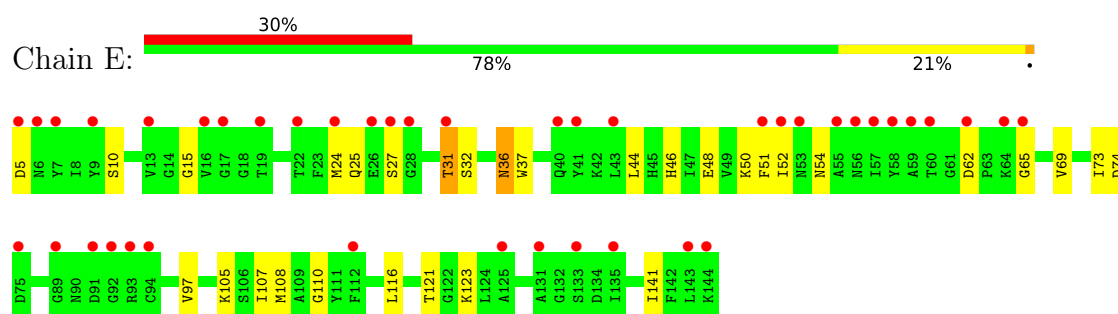
- Molecule 1: lectin



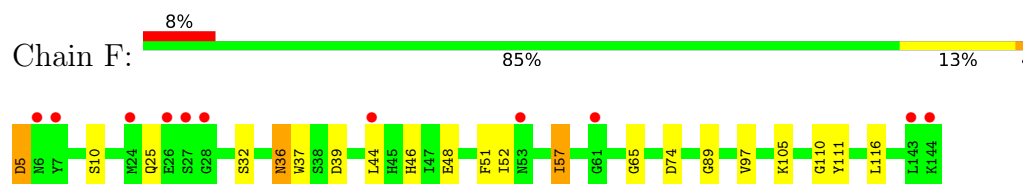
- Molecule 1: lectin



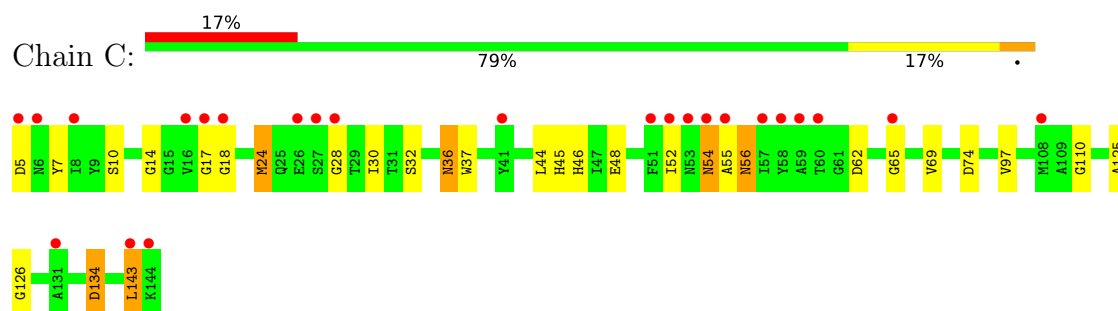
- Molecule 1: lectin



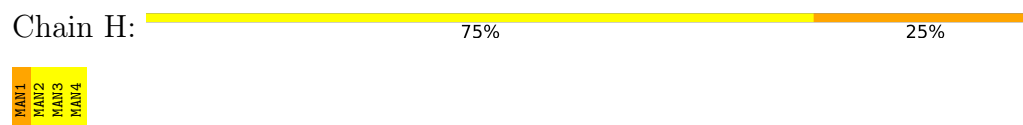
- Molecule 1: lectin



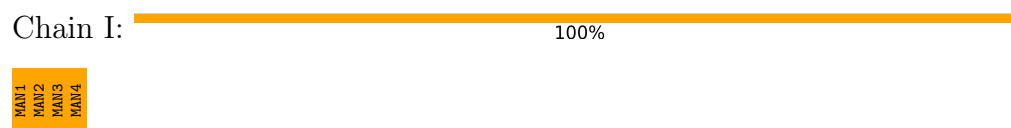
- Molecule 1: lectin



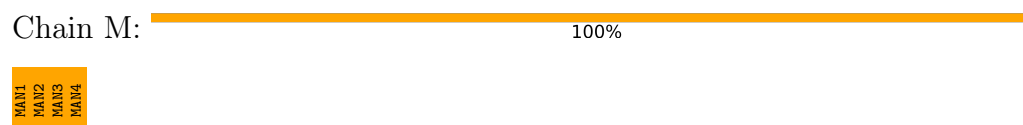
- Molecule 2: alpha-D-mannopyranose-(1-6)-alpha-D-mannopyranose-(1-6)-[alpha-D-mannopyranose-(1-3)]alpha-D-mannopyranose



- Molecule 2: alpha-D-mannopyranose-(1-6)-alpha-D-mannopyranose-(1-6)-[alpha-D-mannopyranose-(1-3)]alpha-D-mannopyranose



- Molecule 2: alpha-D-mannopyranose-(1-6)-alpha-D-mannopyranose-(1-6)-[alpha-D-mannopyranose-(1-3)]alpha-D-mannopyranose



- Molecule 3: α -D-mannopyranose-(1-3)-[α -D-mannopyranose-(1-6)] α -D-mannopyranose

Chain J:  33% 67%



- Molecule 4: α -D-mannopyranose-(1-6)- α -D-mannopyranose

Chain K:  100%



- Molecule 4: α -D-mannopyranose-(1-6)- α -D-mannopyranose

Chain L:  50% 50%



4 Data and refinement statistics

Property	Value	Source
Space group	P 2 21 21	Depositor
Cell constants a, b, c, α , β , γ	51.83Å 168.94Å 169.78Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	85.03 – 2.25 85.03 – 2.25	Depositor EDS
% Data completeness (in resolution range)	100.0 (85.03-2.25) 100.0 (85.03-2.25)	Depositor EDS
R_{merge}	0.21	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.59 (at 2.25Å)	Xtriage
Refinement program	REFMAC 5.8.0258	Depositor
R, R_{free}	0.253 , 0.295 0.256 , 0.296	Depositor DCC
R_{free} test set	3567 reflections (4.96%)	wwPDB-VP
Wilson B-factor (Å ²)	29.4	Xtriage
Anisotropy	0.742	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 29.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.41$, $\langle L^2 \rangle = 0.24$	Xtriage
Estimated twinning fraction	0.068 for -h,l,k	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	7802	wwPDB-VP
Average B, all atoms (Å ²)	36.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.17% 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 ⓘ

5.1 Standard geometry ⓘ

Bond lengths and bond angles in the following residue types are not validated in this section: 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	1.27	4/1078 (0.4%)	1.49	4/1454 (0.3%)
1	B	1.33	3/1084 (0.3%)	1.51	5/1462 (0.3%)
1	C	1.24	1/1072 (0.1%)	1.54	7/1448 (0.5%)
1	D	1.22	3/1077 (0.3%)	1.51	5/1453 (0.3%)
1	E	1.21	3/1084 (0.3%)	1.55	6/1461 (0.4%)
1	F	1.26	1/1080 (0.1%)	1.52	4/1457 (0.3%)
1	G	1.23	3/1051 (0.3%)	1.54	4/1418 (0.3%)
All	All	1.25	18/7526 (0.2%)	1.52	35/10153 (0.3%)

All (18) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	127	ILE	C-O	8.12	1.32	1.23
1	C	17	GLY	C-O	7.53	1.31	1.23
1	B	123	LYS	C-O	6.59	1.32	1.24
1	B	63	PRO	C-O	-6.49	1.15	1.24
1	F	89	GLY	C-O	6.27	1.29	1.23
1	E	107	ILE	C-O	-6.18	1.17	1.24
1	A	23	PHE	C-O	6.17	1.31	1.23
1	G	56	ASN	C-O	6.06	1.31	1.24
1	A	122	GLY	C-O	6.04	1.31	1.23
1	G	85	GLY	C-O	5.96	1.30	1.23
1	A	26	GLU	C-O	5.58	1.30	1.23
1	D	88	LYS	C-O	5.49	1.31	1.23
1	A	33	ILE	C-O	5.40	1.29	1.24
1	E	121	THR	C-O	-5.30	1.17	1.23
1	D	45	HIS	CE1-NE2	5.24	1.37	1.32
1	G	45	HIS	CE1-NE2	5.16	1.37	1.32
1	D	20	PRO	C-O	-5.09	1.18	1.23
1	E	15	GLY	C-O	5.06	1.28	1.23

All (35) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	5	ASP	CA-CB-CG	7.37	119.97	112.60
1	E	31	THR	N-CA-CB	7.12	121.12	110.65
1	C	62	ASP	CA-CB-CG	6.83	119.43	112.60
1	E	31	THR	OG1-CB-CG2	6.16	121.61	109.30
1	G	13	VAL	CA-C-N	6.15	133.46	121.41
1	G	13	VAL	C-N-CA	6.15	133.46	121.41
1	B	15	GLY	CA-C-O	-6.08	116.20	122.28
1	C	28	GLY	N-CA-C	5.88	118.08	111.85
1	D	74	ASP	N-CA-C	-5.67	103.00	110.43
1	A	143	LEU	CA-C-O	-5.66	114.96	121.47
1	G	74	ASP	CA-CB-CG	5.55	118.16	112.60
1	E	62	ASP	CA-CB-CG	5.55	118.15	112.60
1	D	5	ASP	CA-CB-CG	5.52	118.12	112.60
1	A	62	ASP	CA-CB-CG	5.47	118.07	112.60
1	E	74	ASP	N-CA-C	-5.45	103.29	110.43
1	C	62	ASP	CB-CA-C	5.38	115.99	109.85
1	A	74	ASP	N-CA-C	-5.38	103.38	110.43
1	F	74	ASP	N-CA-C	-5.36	103.41	110.43
1	E	51	PHE	CA-C-N	5.36	128.52	120.69
1	E	51	PHE	C-N-CA	5.36	128.52	120.69
1	C	143	LEU	CA-C-O	-5.33	114.84	121.50
1	C	74	ASP	N-CA-C	-5.31	103.47	110.43
1	F	51	PHE	CA-C-N	5.27	128.38	120.69
1	F	51	PHE	C-N-CA	5.27	128.38	120.69
1	D	51	PHE	CA-C-N	5.25	127.73	120.49
1	D	51	PHE	C-N-CA	5.25	127.73	120.49
1	B	74	ASP	N-CA-C	-5.24	103.56	110.43
1	A	62	ASP	CB-CA-C	5.17	115.75	109.85
1	B	51	PHE	CA-C-N	5.14	128.20	120.69
1	B	51	PHE	C-N-CA	5.14	128.20	120.69
1	C	134	ASP	CA-CB-CG	5.12	117.72	112.60
1	G	10	SER	CA-C-O	-5.11	116.14	121.55
1	F	5	ASP	CA-CB-CG	5.10	117.70	112.60
1	D	62	ASP	CA-CB-CG	5.06	117.66	112.60
1	C	14	GLY	CA-C-O	-5.02	117.28	122.25

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts ⓘ

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	1059	0	1038	17	0
1	B	1065	0	1044	19	0
1	C	1053	0	1020	22	0
1	D	1058	0	1036	19	0
1	E	1065	0	1051	18	0
1	F	1061	0	1040	10	0
1	G	1034	0	1006	13	0
2	H	45	0	39	6	1
2	I	45	0	38	6	1
2	M	45	0	38	5	0
3	J	34	0	29	4	0
4	K	23	0	21	3	0
4	L	23	0	21	0	0
5	D	11	0	10	2	0
5	G	12	0	12	7	0
6	D	6	0	8	1	0
7	A	27	0	0	0	0
7	B	33	0	0	3	0
7	C	17	0	0	0	0
7	D	15	0	0	0	0
7	E	31	0	0	2	0
7	F	22	0	0	1	0
7	G	18	0	0	0	0
All	All	7802	0	7451	134	1

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 9.

All (134) 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)
5:G:201:MAN:C2	5:G:201:MAN:C3	1.75	1.58
5:G:201:MAN:C3	5:G:201:MAN:O3	1.66	1.42
2:H:1:MAN:H61	2:H:2:MAN:H3	1.31	1.05
1:E:24:MET:HE1	1:E:141:ILE:HG21	1.53	0.90

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:125:ALA:HB2	1:C:143:LEU:CD1	2.03	0.89
4:K:1:MAN:H62	4:K:2:MAN:H5	1.55	0.88
1:A:31:THR:HB	1:A:52:ILE:HD13	1.55	0.86
1:F:36:ASN:HD21	1:F:65:GLY:H	1.24	0.86
1:G:36:ASN:HD21	1:G:65:GLY:H	1.24	0.85
2:H:1:MAN:H61	2:H:2:MAN:C3	2.00	0.84
1:A:36:ASN:HD21	1:A:65:GLY:H	1.25	0.84
1:B:36:ASN:HD21	1:B:65:GLY:H	1.26	0.84
1:E:36:ASN:HD21	1:E:65:GLY:H	1.24	0.83
1:C:36:ASN:HD21	1:C:65:GLY:H	1.27	0.83
5:G:201:MAN:C2	5:G:201:MAN:H3	2.07	0.82
1:D:36:ASN:HD21	1:D:65:GLY:H	1.26	0.82
1:C:125:ALA:HB2	1:C:143:LEU:HD11	1.64	0.80
5:G:201:MAN:C3	5:G:201:MAN:H2	2.05	0.80
2:I:1:MAN:C6	2:I:2:MAN:H5	2.15	0.77
2:H:1:MAN:O5	2:H:2:MAN:H2	1.90	0.71
3:J:1:MAN:C6	3:J:3:MAN:H3	2.20	0.71
4:K:1:MAN:C6	4:K:2:MAN:H5	2.22	0.69
1:C:54:ASN:ND2	1:C:56:ASN:H	1.90	0.69
1:D:98:LYS:HE2	1:D:108:MET:HE2	1.74	0.69
1:C:7:TYR:CZ	1:C:24:MET:HE1	2.28	0.68
5:G:201:MAN:C2	5:G:201:MAN:C4	2.64	0.68
1:D:36:ASN:HD21	1:D:65:GLY:N	1.92	0.68
1:A:36:ASN:HD21	1:A:65:GLY:N	1.91	0.68
1:F:36:ASN:HD21	1:F:65:GLY:N	1.91	0.67
1:D:18:GLY:HA2	5:D:305:MAN:O2	1.94	0.67
1:E:36:ASN:HD21	1:E:65:GLY:N	1.92	0.67
1:G:36:ASN:HD21	1:G:65:GLY:N	1.92	0.66
1:C:54:ASN:HD22	1:C:55:ALA:N	1.91	0.66
1:B:36:ASN:HD21	1:B:65:GLY:N	1.92	0.66
1:C:36:ASN:HD21	1:C:65:GLY:N	1.92	0.65
2:M:1:MAN:H61	2:M:2:MAN:H3	1.77	0.65
1:C:18:GLY:N	2:M:3:MAN:O3	2.23	0.64
1:B:5:ASP:HB3	7:B:326:HOH:O	1.97	0.64
1:C:54:ASN:HD22	1:C:54:ASN:C	2.07	0.63
1:A:13:VAL:HA	6:D:301:GOL:H31	1.80	0.63
2:I:1:MAN:H61	2:I:2:MAN:H5	1.84	0.60
1:G:132:GLY:HA3	5:G:201:MAN:O1	2.01	0.60
1:B:136:ASP:OD2	2:H:3:MAN:O4	2.17	0.60
3:J:1:MAN:H61	3:J:3:MAN:H3	1.83	0.60
1:D:30:ILE:HD11	1:D:126:GLY:HA2	1.84	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:I:1:MAN:H61	2:I:2:MAN:H3	1.83	0.59
1:B:5:ASP:HB2	7:B:331:HOH:O	2.02	0.59
1:G:88:LYS:HG2	1:G:114:GLU:OE2	2.02	0.59
1:E:25:GLN:HE21	1:E:54:ASN:ND2	2.03	0.57
1:E:108:MET:HE2	7:E:329:HOH:O	2.05	0.56
1:D:77:GLU:OE2	1:D:105:LYS:HD3	2.06	0.55
1:D:98:LYS:CE	1:D:108:MET:HE2	2.35	0.55
1:F:57:ILE:HD12	7:F:312:HOH:O	2.06	0.55
3:J:1:MAN:C6	3:J:3:MAN:C3	2.85	0.55
1:C:30:ILE:HD11	1:C:126:GLY:HA2	1.89	0.54
1:E:31:THR:HG22	1:E:73:ILE:O	2.08	0.54
1:B:18:GLY:HA2	2:H:3:MAN:O2	2.08	0.53
1:D:7:TYR:CE1	1:D:24:MET:HE1	2.44	0.53
1:A:31:THR:HB	1:A:52:ILE:CD1	2.33	0.53
1:E:24:MET:HE1	1:E:141:ILE:CG2	2.33	0.52
1:C:44:LEU:HD11	1:C:97:VAL:HG23	1.93	0.51
1:B:37:TRP:CD1	1:B:110:GLY:HA2	2.46	0.51
1:C:18:GLY:HA2	2:M:3:MAN:O2	2.11	0.51
3:J:1:MAN:H61	3:J:3:MAN:C3	2.40	0.51
1:C:7:TYR:CE1	1:C:24:MET:HE1	2.46	0.51
1:E:37:TRP:CD1	1:E:110:GLY:HA2	2.45	0.51
1:G:44:LEU:HD11	1:G:97:VAL:HG23	1.92	0.51
5:G:201:MAN:C3	5:G:201:MAN:HO3	2.09	0.51
4:K:1:MAN:H62	4:K:2:MAN:C5	2.34	0.51
1:G:16:VAL:HG22	1:G:16:VAL:O	2.11	0.50
1:G:37:TRP:CD1	1:G:110:GLY:HA2	2.46	0.50
1:D:7:TYR:CZ	1:D:24:MET:HE1	2.47	0.49
1:B:128:LYS:HE2	7:B:309:HOH:O	2.11	0.49
1:B:44:LEU:HD11	1:B:97:VAL:HG23	1.94	0.49
1:A:44:LEU:HD11	1:A:97:VAL:HG23	1.95	0.49
1:A:136:ASP:OD1	2:I:3:MAN:O6	2.26	0.49
1:D:37:TRP:CD1	1:D:110:GLY:HA2	2.47	0.48
1:C:54:ASN:HD21	1:C:56:ASN:HB2	1.78	0.48
1:F:37:TRP:CD1	1:F:110:GLY:HA2	2.48	0.48
1:C:37:TRP:CD1	1:C:110:GLY:HA2	2.47	0.48
1:A:37:TRP:CD1	1:A:110:GLY:HA2	2.48	0.48
1:D:44:LEU:HD11	1:D:97:VAL:HG23	1.95	0.48
1:C:45:HIS:HA	1:C:134:ASP:HA	1.95	0.47
1:A:46:HIS:HE1	1:A:48:GLU:OE2	1.98	0.47
1:E:44:LEU:HD11	1:E:97:VAL:HG23	1.96	0.47
1:A:69:VAL:HG12	1:F:116:LEU:HB2	1.97	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:54:ASN:ND2	1:C:54:ASN:C	2.69	0.47
1:E:25:GLN:HA	1:E:25:GLN:OE1	2.14	0.47
1:E:46:HIS:HE1	1:E:48:GLU:OE2	1.97	0.46
1:F:44:LEU:HD11	1:F:97:VAL:HG23	1.98	0.46
1:D:46:HIS:HE1	1:D:48:GLU:OE2	1.99	0.46
1:B:46:HIS:HE1	1:B:48:GLU:OE2	1.99	0.46
1:E:50:LYS:HB3	1:E:50:LYS:HE2	1.71	0.45
2:H:1:MAN:C5	2:H:2:MAN:H2	2.42	0.45
1:C:46:HIS:HE1	1:C:48:GLU:OE2	1.99	0.45
1:F:46:HIS:HE1	1:F:48:GLU:OE2	1.99	0.44
1:A:36:ASN:HD21	1:A:65:GLY:CA	2.31	0.44
1:A:36:ASN:HB3	1:A:46:HIS:HB3	2.00	0.44
1:G:116:LEU:O	1:C:69:VAL:HA	2.17	0.44
1:D:18:GLY:HA2	5:D:305:MAN:HO2	1.80	0.43
1:B:116:LEU:O	1:G:69:VAL:HA	2.19	0.43
1:G:46:HIS:HE1	1:G:48:GLU:OE2	2.00	0.43
1:A:24:MET:HE2	1:A:141:ILE:HD13	2.00	0.43
1:C:143:LEU:CD1	1:C:143:LEU:N	2.81	0.43
1:F:39:ASP:OD1	1:F:111:TYR:OH	2.32	0.43
1:B:36:ASN:HB3	1:B:46:HIS:HB3	2.01	0.43
1:G:36:ASN:HD21	1:G:65:GLY:CA	2.32	0.43
1:C:36:ASN:HD21	1:C:65:GLY:CA	2.32	0.42
1:B:52:ILE:CG2	1:B:53:ASN:HD22	2.32	0.42
1:G:13:VAL:O	1:G:137:ARG:HA	2.19	0.42
1:D:50:LYS:HB3	1:D:50:LYS:HE3	1.68	0.42
2:M:1:MAN:H2	2:M:4:MAN:H5	2.01	0.42
1:G:41:TYR:O	1:G:43:LEU:HG	2.20	0.42
2:M:1:MAN:H61	2:M:2:MAN:C3	2.49	0.42
1:B:52:ILE:HG23	1:B:53:ASN:ND2	2.35	0.42
1:A:36:ASN:ND2	1:A:65:GLY:H	2.05	0.42
1:F:36:ASN:HD21	1:F:65:GLY:CA	2.33	0.42
1:B:36:ASN:HD21	1:B:65:GLY:CA	2.32	0.41
1:A:116:LEU:O	1:D:69:VAL:HA	2.20	0.41
1:B:69:VAL:HG12	1:E:116:LEU:HB2	2.01	0.41
1:E:50:LYS:NZ	7:E:303:HOH:O	2.53	0.41
1:A:69:VAL:HA	1:F:116:LEU:O	2.21	0.41
2:I:1:MAN:C6	2:I:2:MAN:C5	2.95	0.41
1:B:69:VAL:HA	1:E:116:LEU:O	2.21	0.41
1:D:128:LYS:HG2	1:D:141:ILE:HD11	2.02	0.41
1:A:18:GLY:HA2	2:I:3:MAN:O2	2.19	0.41
1:D:116:LEU:HB2	1:E:69:VAL:HG12	2.03	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:54:ASN:HD22	1:C:56:ASN:H	1.64	0.41
1:B:48:GLU:HB3	1:B:57:ILE:HG23	2.03	0.40
1:E:36:ASN:HD21	1:E:65:GLY:CA	2.35	0.40
1:D:36:ASN:HD21	1:D:65:GLY:CA	2.33	0.40
1:B:19:THR:HA	1:B:20:PRO:HD3	1.92	0.40
1:D:36:ASN:HB3	1:D:46:HIS:HB3	2.03	0.40
1:E:123:LYS:HA	1:E:123:LYS:HD2	1.97	0.40

All (1) 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)
2:H:2:MAN:O4	2:I:4:MAN:O3[3_455]	1.65	0.55

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	138/140 (99%)	130 (94%)	7 (5%)	1 (1%)	18	17
1	B	138/140 (99%)	130 (94%)	8 (6%)	0	100	100
1	C	138/140 (99%)	132 (96%)	6 (4%)	0	100	100
1	D	138/140 (99%)	131 (95%)	7 (5%)	0	100	100
1	E	138/140 (99%)	129 (94%)	8 (6%)	1 (1%)	18	17
1	F	138/140 (99%)	130 (94%)	8 (6%)	0	100	100
1	G	135/140 (96%)	124 (92%)	10 (7%)	1 (1%)	18	17
All	All	963/980 (98%)	906 (94%)	54 (6%)	3 (0%)	36	40

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	G	27	SER
1	A	54	ASN
1	E	27	SER

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	111/115 (96%)	106 (96%)	5 (4%)	24	29
1	B	113/115 (98%)	108 (96%)	5 (4%)	25	29
1	C	109/115 (95%)	101 (93%)	8 (7%)	13	12
1	D	111/115 (96%)	106 (96%)	5 (4%)	24	29
1	E	113/115 (98%)	107 (95%)	6 (5%)	20	22
1	F	112/115 (97%)	104 (93%)	8 (7%)	13	12
1	G	108/115 (94%)	98 (91%)	10 (9%)	8	6
All	All	777/805 (96%)	730 (94%)	47 (6%)	17	18

All (47) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	32	SER
1	B	36	ASN
1	B	49	VAL
1	B	52	ILE
1	B	88	LYS
1	G	5	ASP
1	G	10	SER
1	G	25	GLN
1	G	32	SER
1	G	36	ASN
1	G	50	LYS
1	G	52	ILE
1	G	105	LYS
1	G	143	LEU
1	G	144	LYS

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Mol	Chain	Res	Type
1	A	36	ASN
1	A	49	VAL
1	A	52	ILE
1	A	54	ASN
1	A	105	LYS
1	D	5	ASP
1	D	32	SER
1	D	36	ASN
1	D	50	LYS
1	D	52	ILE
1	E	5	ASP
1	E	10	SER
1	E	32	SER
1	E	36	ASN
1	E	52	ILE
1	E	105	LYS
1	F	5	ASP
1	F	10	SER
1	F	25	GLN
1	F	32	SER
1	F	36	ASN
1	F	52	ILE
1	F	57	ILE
1	F	105	LYS
1	C	5	ASP
1	C	10	SER
1	C	24	MET
1	C	32	SER
1	C	36	ASN
1	C	52	ILE
1	C	54	ASN
1	C	56	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (28) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	36	ASN
1	B	46	HIS
1	B	53	ASN
1	B	56	ASN
1	G	36	ASN
1	G	53	ASN

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Mol	Chain	Res	Type
1	G	56	ASN
1	A	36	ASN
1	A	46	HIS
1	D	36	ASN
1	D	46	HIS
1	D	53	ASN
1	D	56	ASN
1	E	25	GLN
1	E	36	ASN
1	E	46	HIS
1	E	53	ASN
1	E	56	ASN
1	F	25	GLN
1	F	36	ASN
1	F	46	HIS
1	F	53	ASN
1	F	56	ASN
1	C	36	ASN
1	C	46	HIS
1	C	53	ASN
1	C	54	ASN
1	C	56	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 ⓘ

19 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	H	1	2	12,12,12	0.51	0	17,17,17	0.80	1 (5%)
2	MAN	H	2	2	11,11,12	0.55	0	15,15,17	0.89	0
2	MAN	H	3	2	11,11,12	0.29	0	15,15,17	0.78	0
2	MAN	H	4	2	11,11,12	0.28	0	15,15,17	0.80	1 (6%)
2	MAN	I	1	2	12,12,12	2.55	5 (41%)	17,17,17	1.92	4 (23%)
2	MAN	I	2	2	11,11,12	2.05	4 (36%)	15,15,17	2.69	10 (66%)
2	MAN	I	3	2	11,11,12	2.41	5 (45%)	15,15,17	0.76	0
2	MAN	I	4	2	11,11,12	3.10	4 (36%)	15,15,17	1.78	4 (26%)
3	MAN	J	1	3	12,12,12	1.36	2 (16%)	17,17,17	2.70	6 (35%)
3	MAN	J	2	3	11,11,12	2.01	4 (36%)	15,15,17	2.02	3 (20%)
3	MAN	J	3	3	11,11,12	3.55	6 (54%)	15,15,17	2.73	8 (53%)
4	MAN	K	1	4	12,12,12	2.84	7 (58%)	17,17,17	2.25	6 (35%)
4	MAN	K	2	4	11,11,12	4.56	7 (63%)	15,15,17	2.82	7 (46%)
4	MAN	L	1	4	12,12,12	0.50	0	17,17,17	0.80	1 (5%)
4	MAN	L	2	4	11,11,12	0.29	0	15,15,17	0.76	0
2	MAN	M	1	2	12,12,12	1.94	4 (33%)	17,17,17	2.48	7 (41%)
2	MAN	M	2	2	11,11,12	3.01	5 (45%)	15,15,17	3.40	8 (53%)
2	MAN	M	3	2	11,11,12	2.72	6 (54%)	15,15,17	2.23	2 (13%)
2	MAN	M	4	2	11,11,12	3.17	4 (36%)	15,15,17	1.94	6 (40%)

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	H	1	2	-	2/2/22/22	0/1/1/1
2	MAN	H	2	2	-	2/2/19/22	0/1/1/1
2	MAN	H	3	2	-	2/2/19/22	0/1/1/1
2	MAN	H	4	2	-	0/2/19/22	0/1/1/1
2	MAN	I	1	2	-	1/2/22/22	0/1/1/1
2	MAN	I	2	2	-	1/2/19/22	0/1/1/1
2	MAN	I	3	2	-	0/2/19/22	0/1/1/1
2	MAN	I	4	2	-	2/2/19/22	0/1/1/1
3	MAN	J	1	3	-	0/2/22/22	0/1/1/1
3	MAN	J	2	3	-	2/2/19/22	0/1/1/1
3	MAN	J	3	3	-	0/2/19/22	0/1/1/1
4	MAN	K	1	4	-	0/2/22/22	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	MAN	K	2	4	-	1/2/19/22	0/1/1/1
4	MAN	L	1	4	-	0/2/22/22	0/1/1/1
4	MAN	L	2	4	-	0/2/19/22	0/1/1/1
2	MAN	M	1	2	-	2/2/22/22	0/1/1/1
2	MAN	M	2	2	-	0/2/19/22	0/1/1/1
2	MAN	M	3	2	-	0/2/19/22	0/1/1/1
2	MAN	M	4	2	-	0/2/19/22	0/1/1/1

All (63) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	K	2	MAN	C2-C3	10.62	1.68	1.52
3	J	3	MAN	C2-C3	8.38	1.65	1.52
2	M	4	MAN	C2-C3	8.37	1.65	1.52
2	I	4	MAN	C2-C3	8.34	1.65	1.52
4	K	2	MAN	O5-C5	7.96	1.58	1.43
2	M	2	MAN	C2-C3	6.40	1.62	1.52
4	K	1	MAN	C1-C2	6.30	1.66	1.52
2	I	1	MAN	C3-C2	5.99	1.67	1.52
2	I	3	MAN	O5-C5	5.03	1.53	1.43
3	J	3	MAN	O5-C5	4.74	1.52	1.43
2	M	2	MAN	O3-C3	4.69	1.54	1.43
2	M	2	MAN	O5-C5	4.56	1.52	1.43
2	I	2	MAN	C2-C3	4.29	1.59	1.52
2	M	3	MAN	O5-C1	4.25	1.50	1.43
4	K	2	MAN	C4-C3	4.06	1.62	1.52
3	J	3	MAN	O3-C3	4.05	1.53	1.43
2	M	3	MAN	O2-C2	-4.01	1.34	1.43
2	M	1	MAN	O5-C1	3.91	1.52	1.42
2	M	4	MAN	O5-C5	3.89	1.51	1.43
2	I	1	MAN	O5-C5	3.88	1.53	1.44
3	J	2	MAN	O5-C5	3.87	1.51	1.43
2	M	3	MAN	O4-C4	3.81	1.52	1.43
2	M	3	MAN	C1-C2	-3.72	1.43	1.52
2	I	3	MAN	O4-C4	3.65	1.52	1.43
4	K	1	MAN	C3-C2	3.62	1.61	1.52
2	I	4	MAN	O4-C4	-3.51	1.34	1.43
3	J	2	MAN	O5-C1	3.38	1.49	1.43
4	K	2	MAN	O3-C3	3.33	1.51	1.43
2	M	3	MAN	C4-C3	3.31	1.60	1.52
2	M	1	MAN	C4-C5	-3.24	1.46	1.53
3	J	3	MAN	C4-C5	3.23	1.59	1.53

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	I	2	MAN	C1-C2	3.21	1.59	1.52
2	M	4	MAN	O3-C3	3.09	1.50	1.43
2	I	4	MAN	O5-C1	2.98	1.48	1.43
4	K	2	MAN	O5-C1	2.92	1.48	1.43
2	I	1	MAN	O2-C2	2.89	1.50	1.43
2	I	1	MAN	C4-C5	-2.83	1.47	1.53
2	M	4	MAN	O2-C2	2.82	1.49	1.43
4	K	2	MAN	O4-C4	2.79	1.49	1.43
4	K	1	MAN	C4-C3	2.75	1.59	1.52
3	J	3	MAN	O2-C2	2.74	1.49	1.43
2	I	3	MAN	C2-C3	-2.69	1.48	1.52
4	K	1	MAN	O2-C2	2.66	1.49	1.43
3	J	2	MAN	C2-C3	2.63	1.56	1.52
2	M	1	MAN	O4-C4	2.63	1.49	1.43
4	K	1	MAN	O1-C1	2.63	1.47	1.39
4	K	1	MAN	O3-C3	2.61	1.49	1.43
2	I	1	MAN	O4-C4	2.59	1.49	1.43
3	J	2	MAN	O4-C4	2.53	1.49	1.43
3	J	1	MAN	O1-C1	2.46	1.47	1.39
4	K	1	MAN	O5-C1	2.44	1.48	1.42
2	M	2	MAN	O5-C1	-2.42	1.39	1.43
2	I	3	MAN	C1-C2	2.36	1.57	1.52
2	I	3	MAN	O6-C6	2.31	1.52	1.42
3	J	1	MAN	C4-C3	2.28	1.58	1.52
2	M	1	MAN	C4-C3	2.22	1.58	1.52
2	I	2	MAN	C4-C3	2.20	1.58	1.52
2	M	3	MAN	C2-C3	2.19	1.55	1.52
2	I	4	MAN	O2-C2	2.18	1.47	1.43
2	M	2	MAN	C6-C5	-2.11	1.44	1.51
4	K	2	MAN	C6-C5	2.08	1.58	1.51
2	I	2	MAN	O5-C1	2.03	1.47	1.43
3	J	3	MAN	C1-C2	2.02	1.57	1.52

All (74) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	M	2	MAN	O5-C5-C6	-6.72	94.59	107.66
2	M	3	MAN	O3-C3-C2	-6.61	96.57	110.05
4	K	2	MAN	O5-C5-C6	6.57	120.44	107.66
3	J	1	MAN	C6-C5-C4	-5.86	98.64	113.02
4	K	1	MAN	O5-C5-C6	5.65	120.45	106.44
3	J	2	MAN	C1-O5-C5	-5.35	105.02	112.19

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	J	1	MAN	O5-C5-C4	5.31	119.27	109.70
4	K	2	MAN	O2-C2-C3	5.14	120.79	110.15
2	M	1	MAN	C3-C4-C5	-5.06	101.06	110.23
2	M	2	MAN	O2-C2-C1	-4.93	97.93	109.22
3	J	1	MAN	O5-C5-C6	-4.86	94.41	106.44
2	I	2	MAN	O6-C6-C5	-4.79	95.02	111.33
2	M	2	MAN	C1-C2-C3	4.71	116.50	109.64
2	I	2	MAN	C1-O5-C5	-4.70	105.89	112.19
3	J	3	MAN	O3-C3-C2	4.63	119.50	110.05
2	M	2	MAN	O3-C3-C2	4.60	119.44	110.05
2	M	1	MAN	O4-C4-C3	4.54	121.08	110.38
4	K	2	MAN	O3-C3-C2	4.25	118.73	110.05
3	J	3	MAN	O5-C5-C6	-4.24	99.41	107.66
2	M	2	MAN	C1-O5-C5	-4.01	106.81	112.19
3	J	2	MAN	O6-C6-C5	-3.99	97.76	111.33
3	J	3	MAN	C2-C3-C4	-3.98	103.86	110.86
3	J	3	MAN	O3-C3-C4	3.92	119.61	110.38
2	I	4	MAN	O2-C2-C3	3.89	118.21	110.15
2	M	1	MAN	C4-C3-C2	-3.82	104.12	110.83
2	I	1	MAN	C3-C4-C5	-3.66	103.60	110.23
2	I	2	MAN	C3-C4-C5	-3.55	103.79	110.23
2	M	2	MAN	C3-C4-C5	-3.55	103.80	110.23
3	J	3	MAN	O6-C6-C5	-3.54	99.28	111.33
2	M	3	MAN	C2-C3-C4	3.51	117.03	110.86
4	K	2	MAN	C3-C4-C5	-3.44	104.00	110.23
2	I	1	MAN	O3-C3-C4	-3.41	102.33	110.38
2	M	1	MAN	O5-C5-C6	3.35	114.75	106.44
4	K	1	MAN	C6-C5-C4	-3.35	104.80	113.02
2	M	4	MAN	O2-C2-C3	3.35	117.08	110.15
2	M	1	MAN	O5-C1-C2	3.31	116.12	110.30
3	J	1	MAN	O4-C4-C3	3.27	118.08	110.38
2	I	4	MAN	O6-C6-C5	-3.26	100.24	111.33
2	M	2	MAN	C2-C3-C4	-3.24	105.17	110.86
2	I	1	MAN	O6-C6-C5	-3.14	100.63	111.33
3	J	3	MAN	C6-C5-C4	-3.11	105.38	113.02
2	M	4	MAN	C6-C5-C4	-3.05	105.52	113.02
4	K	1	MAN	O5-C1-C2	3.03	115.62	110.30
2	I	2	MAN	C2-C3-C4	-2.95	105.67	110.86
4	K	1	MAN	C1-C2-C3	2.93	116.33	110.36
2	M	4	MAN	C3-C4-C5	2.92	115.52	110.23
2	I	1	MAN	O2-C2-C3	2.89	117.18	110.38
2	I	2	MAN	O3-C3-C2	2.83	115.82	110.05

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	I	2	MAN	O4-C4-C3	2.67	116.67	110.38
3	J	1	MAN	O1-C1-O5	-2.65	102.55	110.41
2	M	2	MAN	O6-C6-C5	-2.64	102.36	111.33
2	I	4	MAN	C1-C2-C3	-2.61	105.84	109.64
2	M	1	MAN	C6-C5-C4	-2.51	106.85	113.02
2	M	4	MAN	O3-C3-C2	2.50	115.16	110.05
3	J	3	MAN	O2-C2-C3	2.50	115.33	110.15
2	I	2	MAN	C1-C2-C3	2.48	113.25	109.64
4	K	1	MAN	C3-C4-C5	-2.44	105.81	110.23
2	I	4	MAN	O4-C4-C5	-2.41	103.39	109.32
3	J	1	MAN	C4-C3-C2	-2.41	106.60	110.83
4	K	2	MAN	O5-C1-C2	-2.41	105.05	110.79
3	J	2	MAN	O5-C5-C6	2.32	112.17	107.66
2	I	2	MAN	O2-C2-C1	-2.26	104.05	109.22
4	L	1	MAN	C1-O5-C5	-2.23	109.34	113.65
3	J	3	MAN	C3-C4-C5	-2.21	106.22	110.23
2	M	1	MAN	O5-C5-C4	-2.21	105.72	109.70
4	K	2	MAN	O2-C2-C1	-2.20	104.18	109.22
2	H	1	MAN	C1-O5-C5	-2.20	109.40	113.65
4	K	2	MAN	O3-C3-C4	2.16	115.46	110.38
2	I	2	MAN	O5-C5-C4	-2.15	105.60	110.83
2	H	4	MAN	C1-O5-C5	-2.14	109.32	112.19
4	K	1	MAN	O2-C2-C1	2.13	114.16	109.25
2	M	4	MAN	C1-O5-C5	2.10	115.00	112.19
2	M	4	MAN	C2-C3-C4	2.09	114.54	110.86
2	I	2	MAN	O5-C1-C2	2.01	115.59	110.79

There are no chirality outliers.

All (15) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	J	2	MAN	C4-C5-C6-O6
2	H	2	MAN	O5-C5-C6-O6
2	H	2	MAN	C4-C5-C6-O6
2	H	1	MAN	O5-C5-C6-O6
2	M	1	MAN	O5-C5-C6-O6
2	H	3	MAN	C4-C5-C6-O6
2	I	2	MAN	C4-C5-C6-O6
2	H	3	MAN	O5-C5-C6-O6
2	I	1	MAN	O5-C5-C6-O6
2	I	4	MAN	O5-C5-C6-O6
3	J	2	MAN	O5-C5-C6-O6

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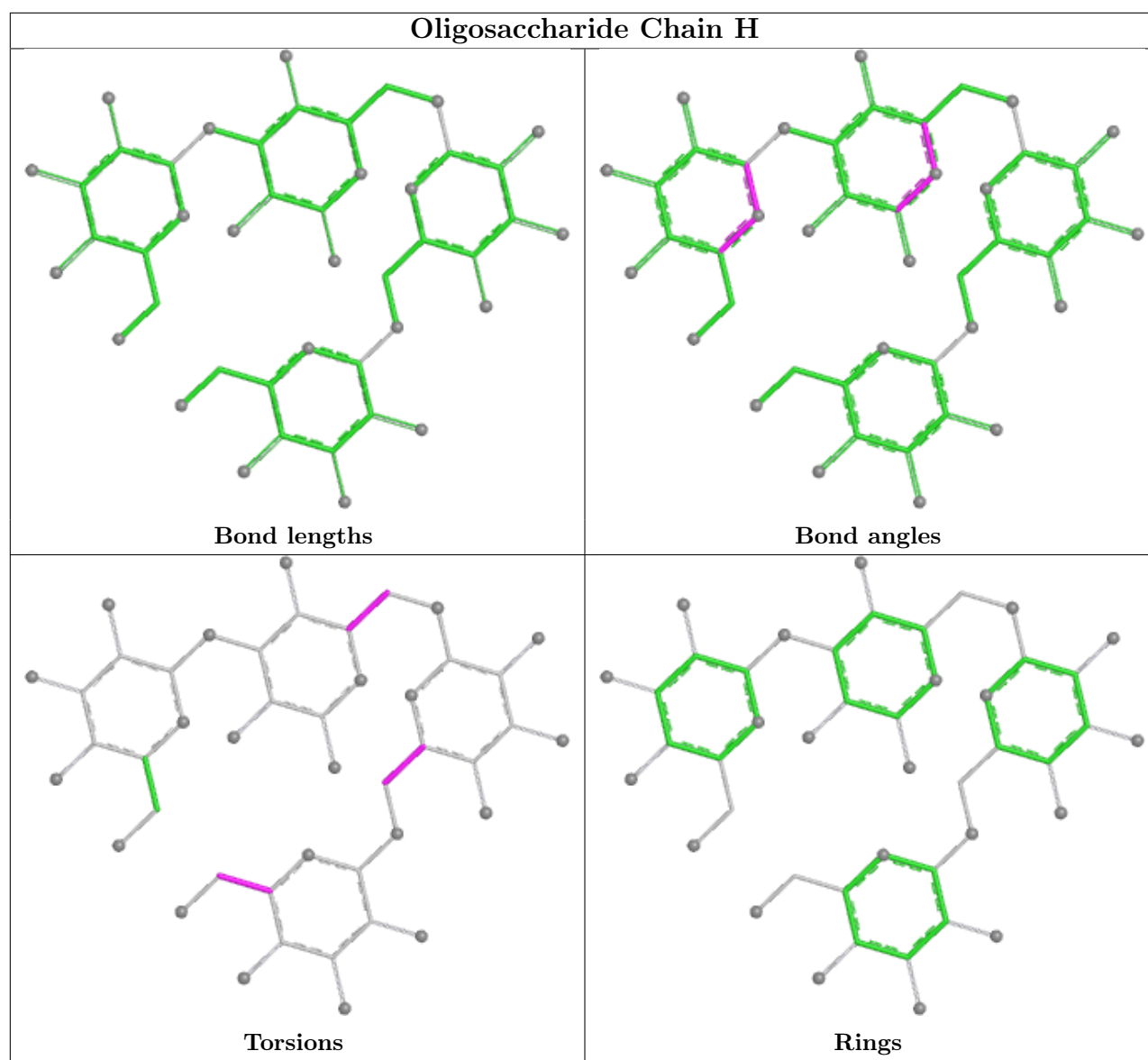
Mol	Chain	Res	Type	Atoms
2	M	1	MAN	C4-C5-C6-O6
2	I	4	MAN	C4-C5-C6-O6
4	K	2	MAN	C4-C5-C6-O6
2	H	1	MAN	C4-C5-C6-O6

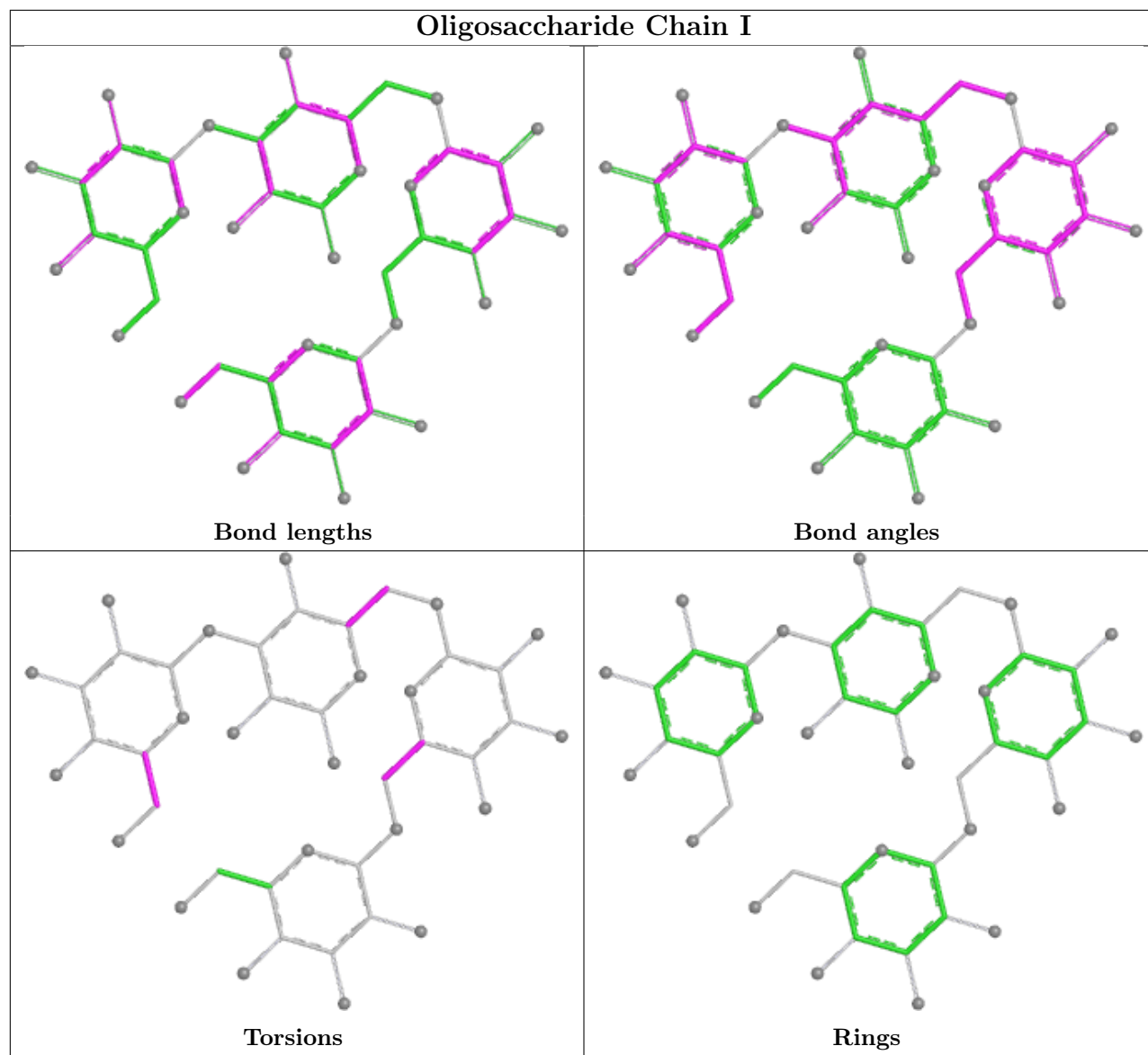
There are no ring outliers.

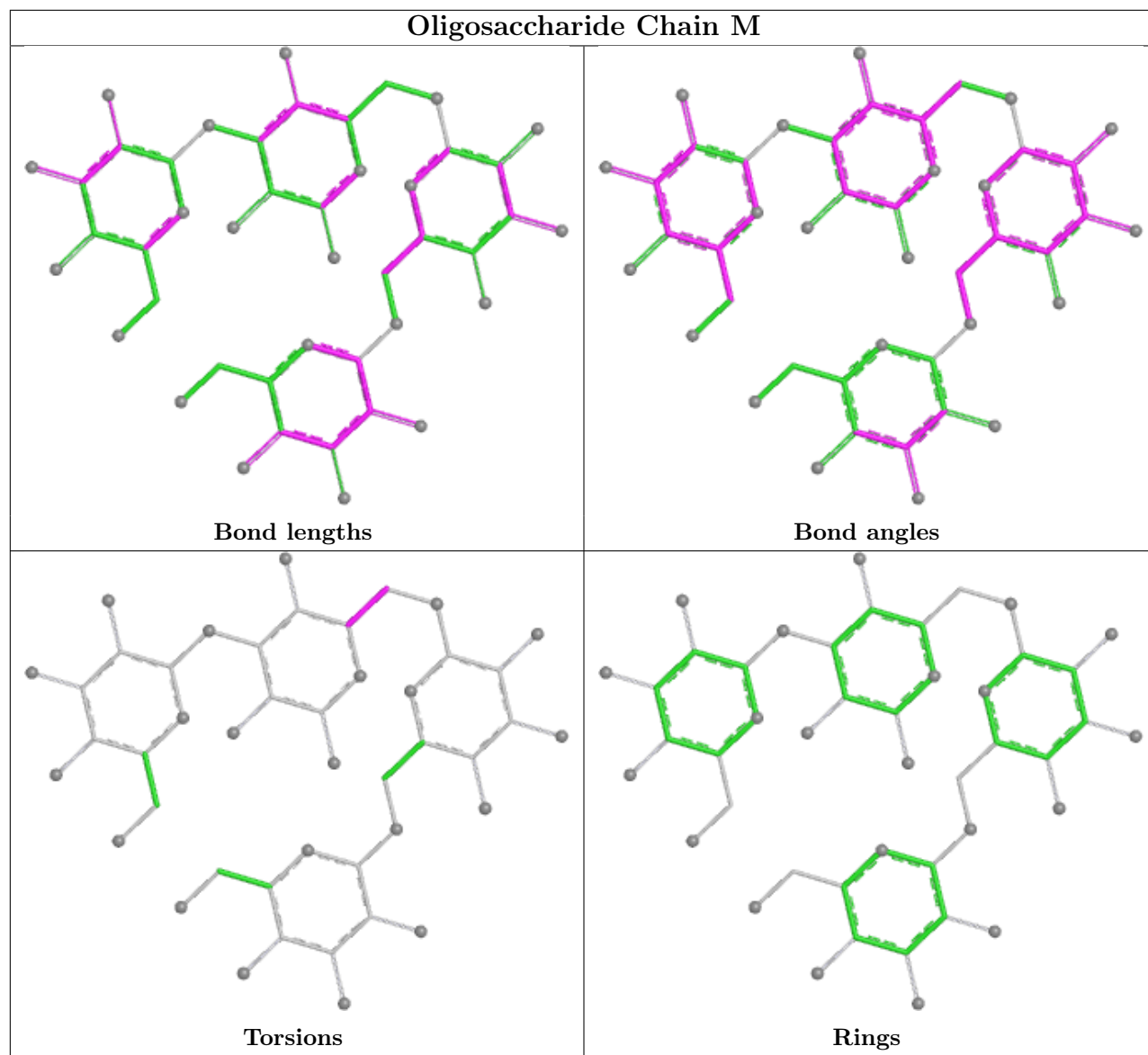
15 monomers are involved in 25 short contacts:

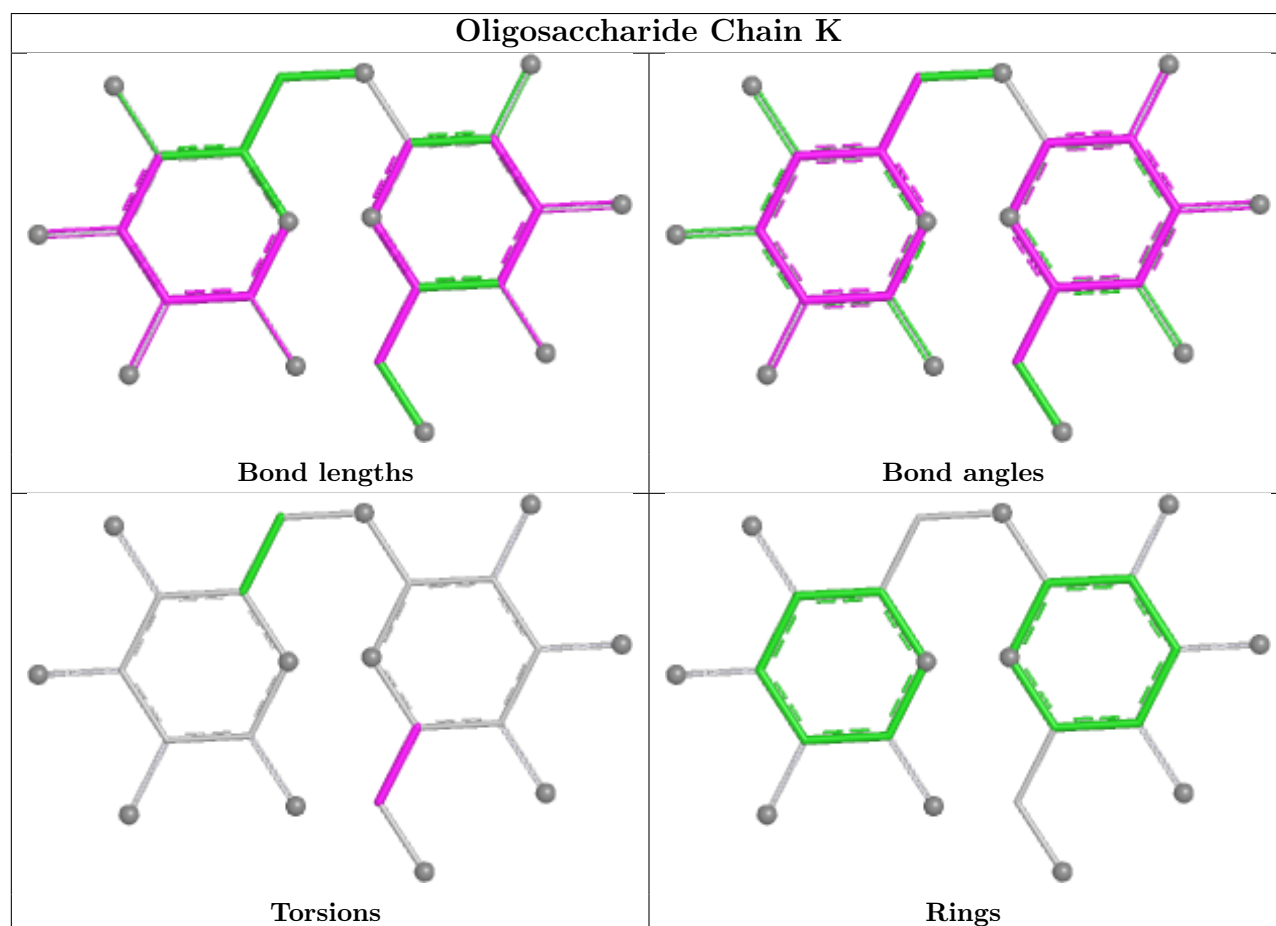
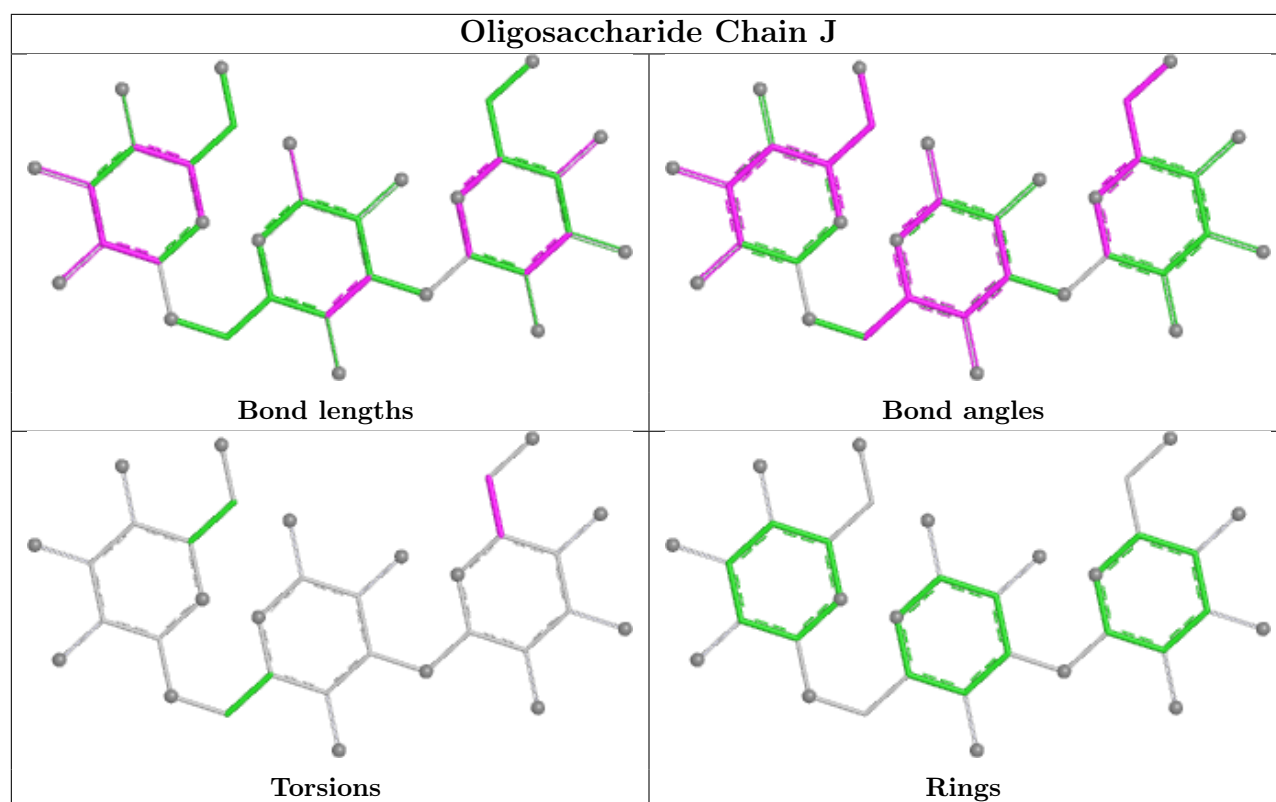
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	I	1	MAN	4	0
2	M	4	MAN	1	0
4	K	2	MAN	3	0
2	M	2	MAN	2	0
2	H	1	MAN	4	0
2	I	2	MAN	4	0
3	J	3	MAN	4	0
2	H	3	MAN	2	0
2	H	2	MAN	4	1
2	I	3	MAN	2	0
2	I	4	MAN	0	1
3	J	1	MAN	4	0
4	K	1	MAN	3	0
2	M	1	MAN	3	0
2	M	3	MAN	2	0

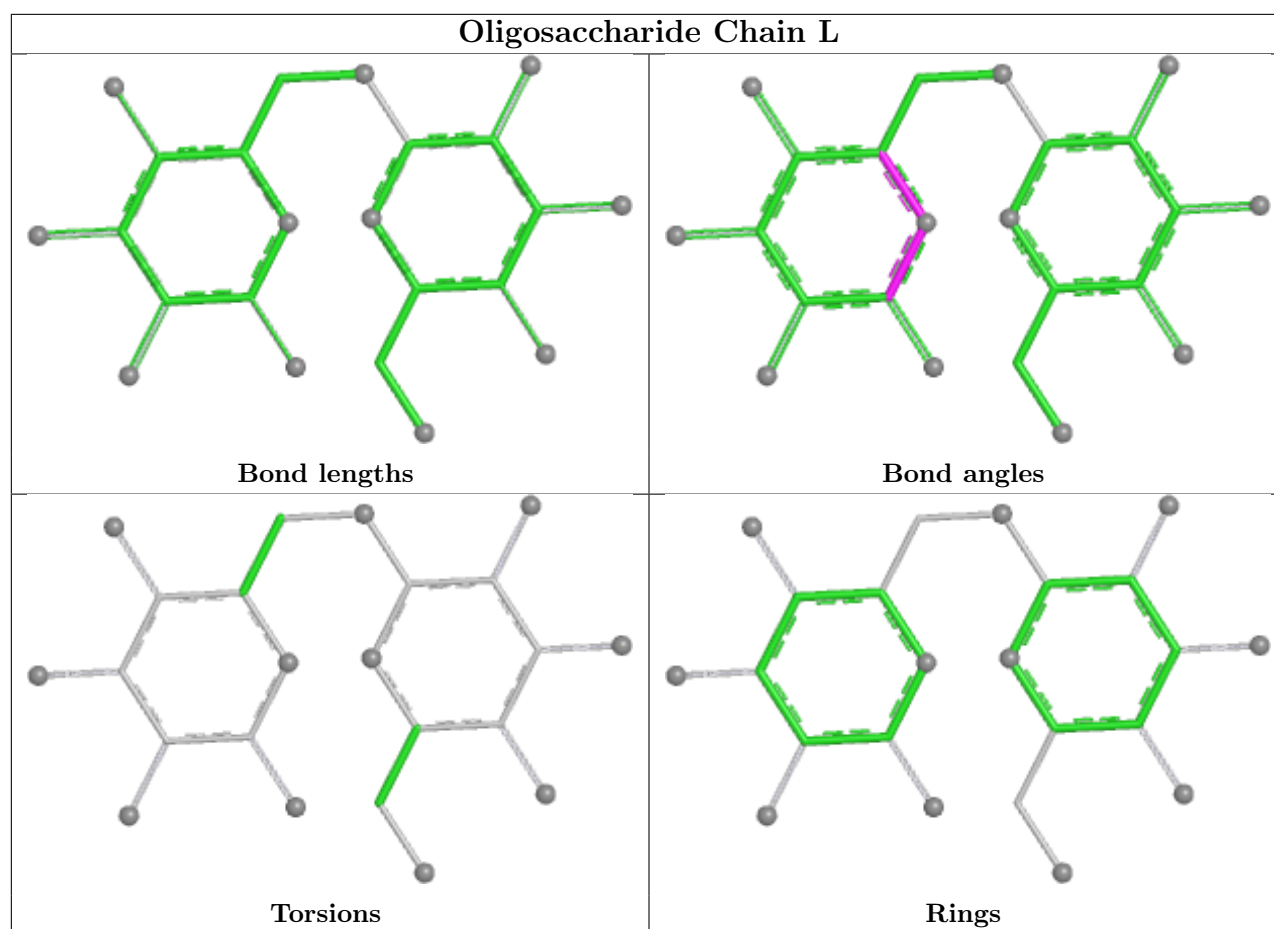
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for oligosaccharide.











5.6 Ligand geometry [i](#)

3 ligands are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
5	MAN	G	201	-	12,12,12	4.46	7 (58%)	17,17,17	4.43	11 (64%)
5	MAN	D	305	-	11,11,12	2.76	4 (36%)	15,15,17	1.91	4 (26%)
6	GOL	D	301	-	5,5,5	0.09	0	5,5,5	0.40	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns.

'-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
5	MAN	G	201	-	-	2/2/22/22	0/1/1/1
5	MAN	D	305	-	-	0/2/19/22	0/1/1/1
6	GOL	D	301	-	-	1/4/4/4	-

All (11) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	G	201	MAN	O3-C3	9.30	1.66	1.43
5	G	201	MAN	C3-C2	8.81	1.75	1.52
5	D	305	MAN	O4-C4	5.32	1.56	1.43
5	D	305	MAN	C4-C3	5.29	1.66	1.52
5	G	201	MAN	C4-C5	5.19	1.64	1.53
5	G	201	MAN	O4-C4	4.87	1.55	1.43
5	D	305	MAN	C4-C5	-3.41	1.45	1.53
5	G	201	MAN	O5-C5	2.46	1.50	1.44
5	G	201	MAN	O1-C1	2.42	1.47	1.39
5	D	305	MAN	C2-C3	2.41	1.56	1.52
5	G	201	MAN	C4-C3	-2.35	1.46	1.52

All (15) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	G	201	MAN	O4-C4-C5	7.40	127.55	109.32
5	G	201	MAN	C3-C4-C5	-7.15	97.27	110.23
5	G	201	MAN	O3-C3-C2	6.76	126.32	110.38
5	G	201	MAN	C1-O5-C5	-6.19	101.68	113.65
5	G	201	MAN	C6-C5-C4	5.93	127.58	113.02
5	G	201	MAN	O4-C4-C3	-5.93	96.41	110.38
5	G	201	MAN	O2-C2-C1	-5.09	97.48	109.25
5	D	305	MAN	O3-C3-C2	-4.68	100.50	110.05
5	G	201	MAN	C1-C2-C3	3.62	117.75	110.36
5	G	201	MAN	O2-C2-C3	3.48	118.58	110.38
5	D	305	MAN	C1-C2-C3	-3.37	104.73	109.64
5	G	201	MAN	O5-C1-C2	-3.36	104.39	110.30
5	G	201	MAN	O1-C1-O5	2.54	117.95	110.41
5	D	305	MAN	O3-C3-C4	2.27	115.73	110.38
5	D	305	MAN	O5-C1-C2	-2.03	105.95	110.79

There are no chirality outliers.

All (3) torsion outliers are listed below:

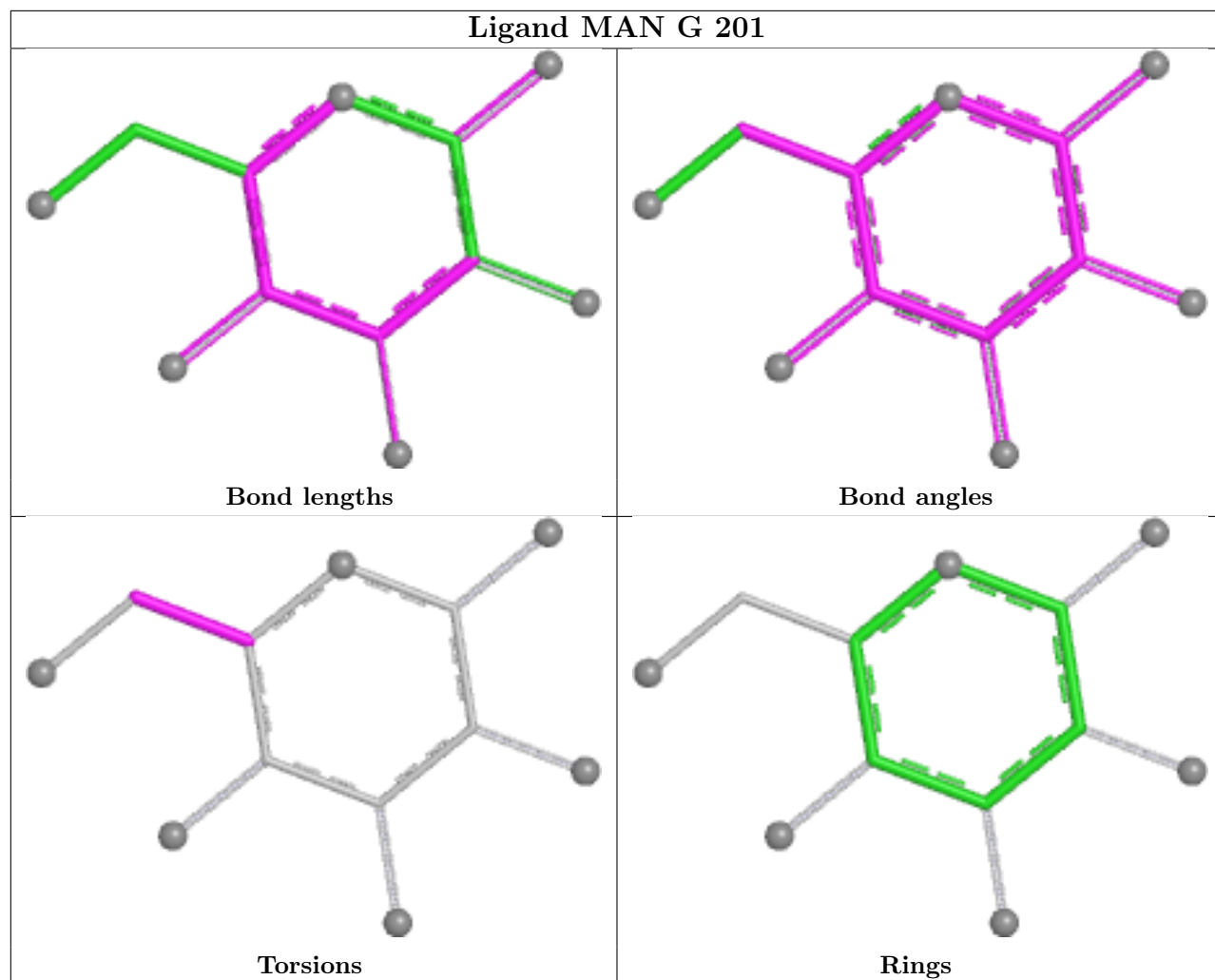
Mol	Chain	Res	Type	Atoms
5	G	201	MAN	O5-C5-C6-O6
5	G	201	MAN	C4-C5-C6-O6
6	D	301	GOL	O1-C1-C2-O2

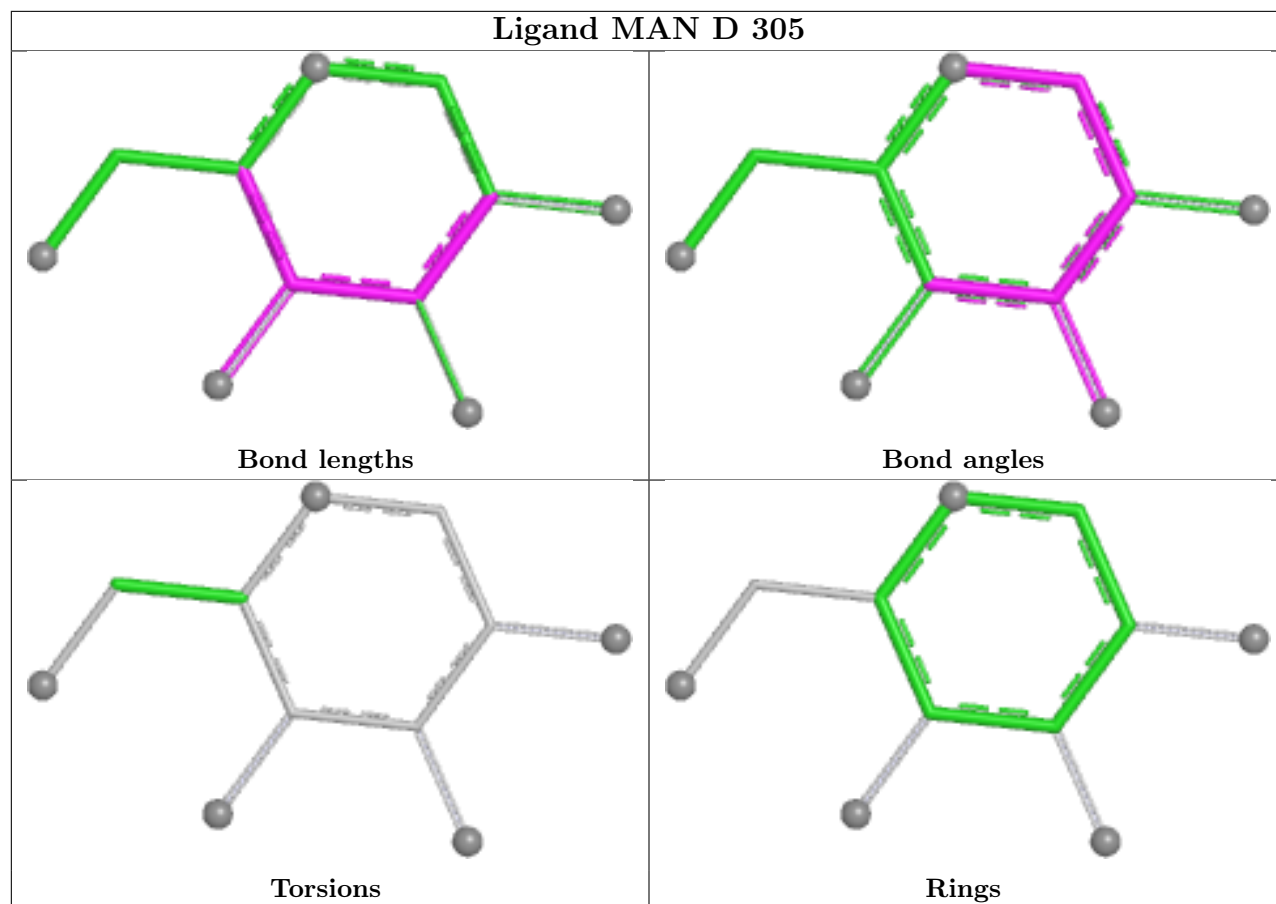
There are no ring outliers.

3 monomers are involved in 10 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	G	201	MAN	7	0
5	D	305	MAN	2	0
6	D	301	GOL	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 all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.





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	140/140 (100%)	-0.17	3 (2%) 63 64	12, 22, 42, 57	0
1	B	140/140 (100%)	0.15	6 (4%) 40 39	13, 25, 48, 69	0
1	C	140/140 (100%)	1.01	24 (17%) 4 3	18, 36, 61, 85	0
1	D	140/140 (100%)	0.55	13 (9%) 14 13	16, 32, 51, 74	0
1	E	140/140 (100%)	1.41	42 (30%) 1 1	22, 42, 72, 109	0
1	F	140/140 (100%)	0.44	11 (7%) 18 17	14, 30, 62, 87	0
1	G	139/140 (99%)	1.71	49 (35%) 1 0	23, 44, 78, 103	0
All	All	979/980 (99%)	0.73	148 (15%) 5 5	12, 33, 65, 109	0

All (148) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	G	27	SER	6.7
1	G	7	TYR	6.6
1	E	26	GLU	6.3
1	E	28	GLY	6.0
1	F	27	SER	5.6
1	G	17	GLY	5.4
1	F	26	GLU	5.2
1	A	53	ASN	5.2
1	G	18	GLY	5.2
1	A	5	ASP	5.1
1	G	16	VAL	4.9
1	G	135	ILE	4.9
1	D	89	GLY	4.8
1	G	14	GLY	4.7
1	G	52	ILE	4.6
1	G	20	PRO	4.5
1	E	27	SER	4.5

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Mol	Chain	Res	Type	RSRZ
1	E	144	LYS	4.3
1	G	94	CYS	4.2
1	D	144	LYS	4.2
1	C	27	SER	4.1
1	G	131	ALA	4.1
1	C	144	LYS	4.1
1	E	52	ILE	4.1
1	G	55	ALA	4.0
1	G	58	TYR	4.0
1	G	144	LYS	3.9
1	G	143	LEU	3.9
1	C	28	GLY	3.9
1	C	6	ASN	3.9
1	E	57	ILE	3.7
1	G	26	GLU	3.7
1	G	23	PHE	3.7
1	B	144	LYS	3.7
1	E	6	ASN	3.6
1	G	6	ASN	3.5
1	G	22	THR	3.5
1	G	137	ARG	3.5
1	E	135	ILE	3.5
1	G	44	LEU	3.5
1	E	89	GLY	3.4
1	E	92	GLY	3.4
1	G	133	SER	3.4
1	B	5	ASP	3.4
1	E	7	TYR	3.3
1	G	9	TYR	3.3
1	E	55	ALA	3.3
1	G	90	ASN	3.3
1	B	53	ASN	3.3
1	G	122	GLY	3.2
1	G	21	PHE	3.2
1	E	16	VAL	3.2
1	G	28	GLY	3.1
1	E	5	ASP	3.1
1	G	53	ASN	3.1
1	D	6	ASN	3.1
1	C	57	ILE	3.1
1	E	94	CYS	3.1
1	F	6	ASN	3.1

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Mol	Chain	Res	Type	RSRZ
1	E	9	TYR	3.1
1	B	52	ILE	3.1
1	G	132	GLY	3.0
1	F	28	GLY	3.0
1	E	64	LYS	3.0
1	C	59	ALA	3.0
1	D	16	VAL	3.0
1	D	27	SER	3.0
1	E	60	THR	3.0
1	G	8	ILE	2.9
1	G	91	ASP	2.9
1	C	55	ALA	2.9
1	F	143	LEU	2.9
1	G	54	ASN	2.9
1	G	15	GLY	2.8
1	E	143	LEU	2.8
1	D	5	ASP	2.8
1	G	41	TYR	2.8
1	E	40	GLN	2.8
1	E	41	TYR	2.8
1	F	144	LYS	2.8
1	E	125	ALA	2.8
1	D	17	GLY	2.8
1	G	60	THR	2.7
1	E	62	ASP	2.7
1	C	65	GLY	2.7
1	C	26	GLU	2.7
1	G	138	LEU	2.7
1	C	58	TYR	2.7
1	G	87	LYS	2.6
1	E	58	TYR	2.6
1	D	90	ASN	2.6
1	C	18	GLY	2.6
1	D	28	GLY	2.6
1	C	8	ILE	2.6
1	D	7	TYR	2.6
1	E	131	ALA	2.6
1	G	19	THR	2.6
1	C	143	LEU	2.5
1	E	91	ASP	2.5
1	E	24	MET	2.5
1	E	53	ASN	2.5

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Mol	Chain	Res	Type	RSRZ
1	G	123	LYS	2.5
1	C	131	ALA	2.5
1	E	75	ASP	2.5
1	C	17	GLY	2.5
1	E	19	THR	2.4
1	E	51	PHE	2.4
1	C	52	ILE	2.4
1	F	24	MET	2.4
1	C	108	MET	2.4
1	D	91	ASP	2.3
1	C	5	ASP	2.3
1	E	31	THR	2.3
1	G	43	LEU	2.3
1	C	53	ASN	2.3
1	D	26	GLU	2.3
1	E	56	ASN	2.3
1	G	59	ALA	2.3
1	E	59	ALA	2.3
1	G	25	GLN	2.3
1	C	51	PHE	2.2
1	E	93	ARG	2.2
1	E	65	GLY	2.2
1	F	7	TYR	2.2
1	B	75	ASP	2.2
1	G	5	ASP	2.2
1	G	56	ASN	2.2
1	G	75	ASP	2.2
1	E	43	LEU	2.2
1	F	44	LEU	2.2
1	A	144	LYS	2.2
1	B	55	ALA	2.2
1	C	41	TYR	2.1
1	D	143	LEU	2.1
1	C	54	ASN	2.1
1	G	125	ALA	2.1
1	G	92	GLY	2.1
1	G	124	LEU	2.1
1	C	16	VAL	2.1
1	E	13	VAL	2.1
1	E	112	PHE	2.1
1	F	61	GLY	2.1
1	E	22	THR	2.0

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Mol	Chain	Res	Type	RSRZ
1	E	133	SER	2.0
1	F	53	ASN	2.0
1	E	17	GLY	2.0
1	C	60	THR	2.0
1	G	88	LYS	2.0

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

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

6.3 Carbohydrates [i](#)

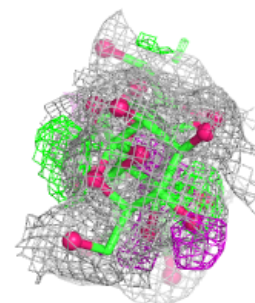
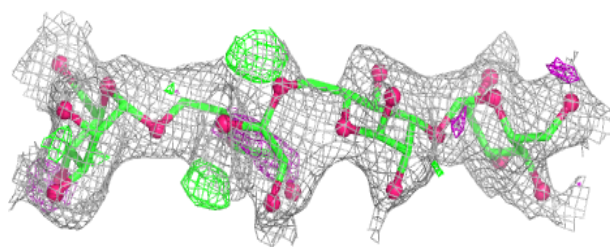
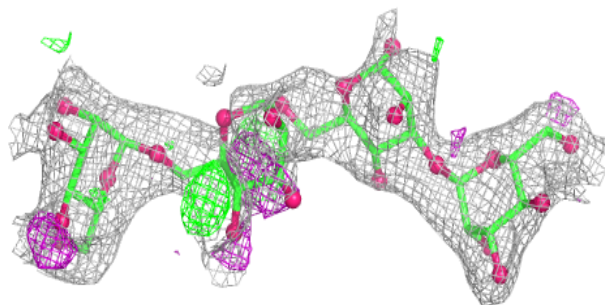
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
4	MAN	L	1	12/12	0.48	0.22	46,78,82,85	0
4	MAN	K	1	12/12	0.62	0.21	61,82,98,98	0
2	MAN	H	2	11/12	0.63	0.23	33,70,85,86	0
2	MAN	H	1	12/12	0.67	0.22	59,69,83,93	0
2	MAN	I	1	12/12	0.73	0.19	66,76,95,99	0
2	MAN	M	1	12/12	0.76	0.17	58,67,75,76	0
2	MAN	I	2	11/12	0.79	0.17	36,72,91,95	0
3	MAN	J	1	12/12	0.81	0.13	58,66,80,89	0
3	MAN	J	3	11/12	0.81	0.15	38,62,73,77	0
4	MAN	K	2	11/12	0.83	0.16	45,63,71,72	0
3	MAN	J	2	11/12	0.85	0.13	54,63,69,69	0
2	MAN	H	4	11/12	0.86	0.14	30,46,57,57	0
2	MAN	I	4	11/12	0.87	0.13	36,44,58,59	0
2	MAN	M	2	11/12	0.87	0.10	43,55,61,63	0
4	MAN	L	2	11/12	0.87	0.12	26,38,45,46	0
2	MAN	H	3	11/12	0.88	0.11	17,21,28,31	0
2	MAN	M	4	11/12	0.92	0.10	31,47,60,67	0
2	MAN	M	3	11/12	0.94	0.09	25,32,38,40	0
2	MAN	I	3	11/12	0.95	0.07	19,22,29,30	0

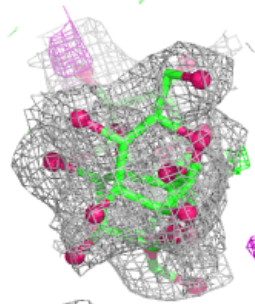
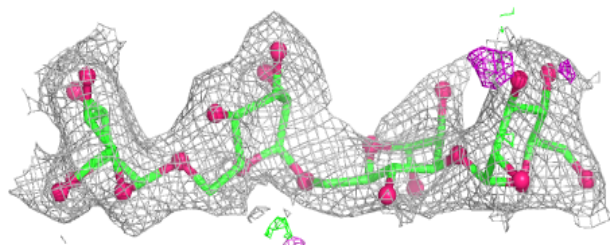
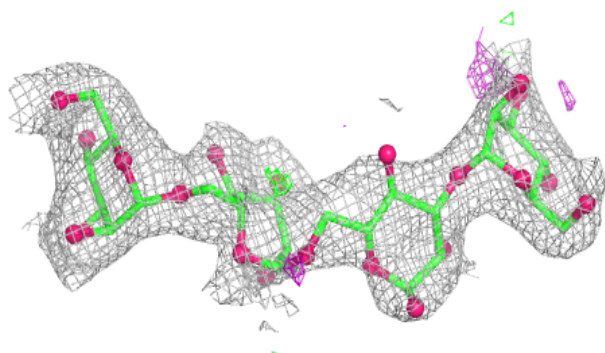
The following is a graphical depiction of the model fit to experimental electron density for oligosaccharide. Each fit is shown from different orientation to approximate a three-dimensional view.

Electron density around Chain H:

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

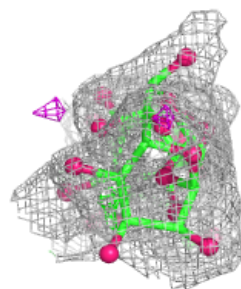
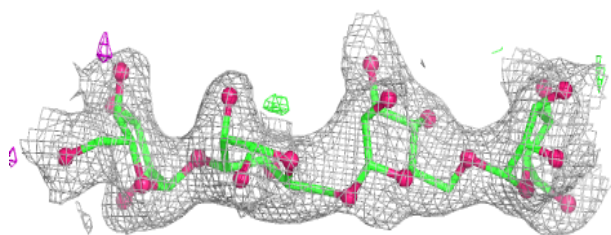
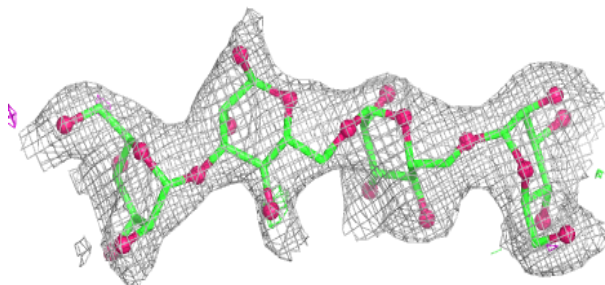
**Electron density around Chain I:**

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

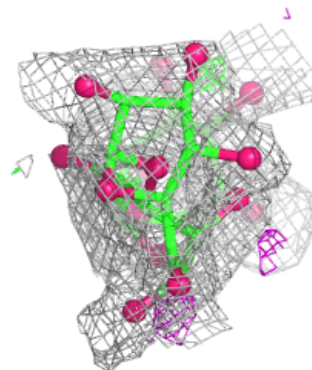
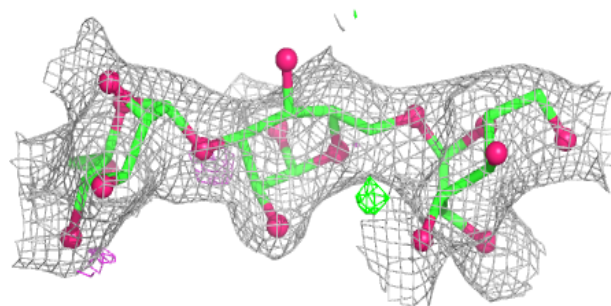
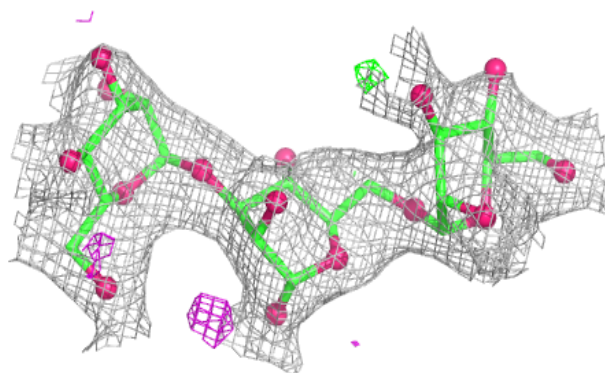


Electron density around Chain M:

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

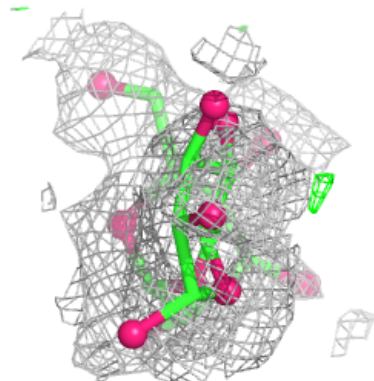
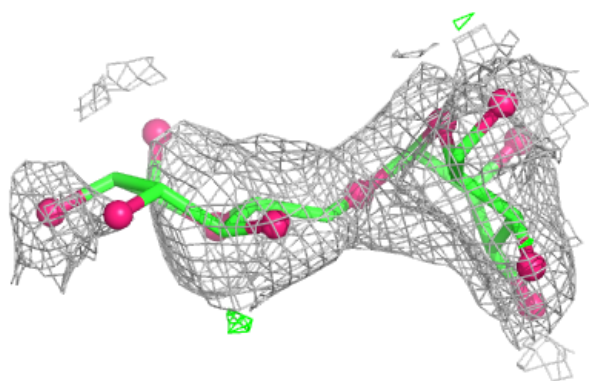
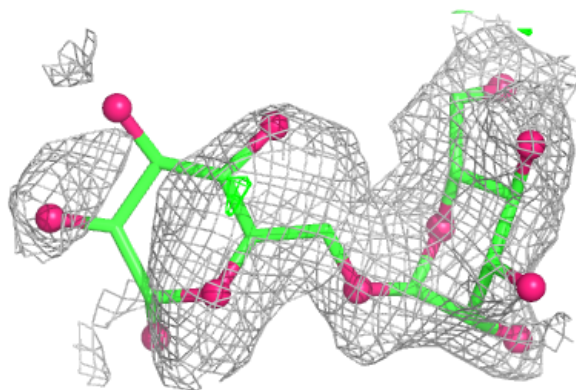
**Electron density around Chain J:**

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

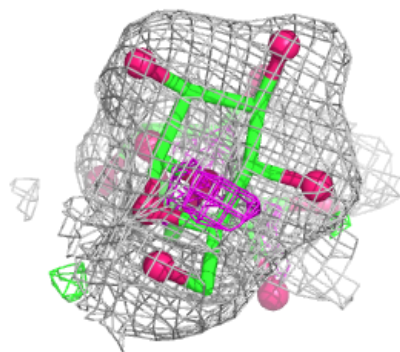
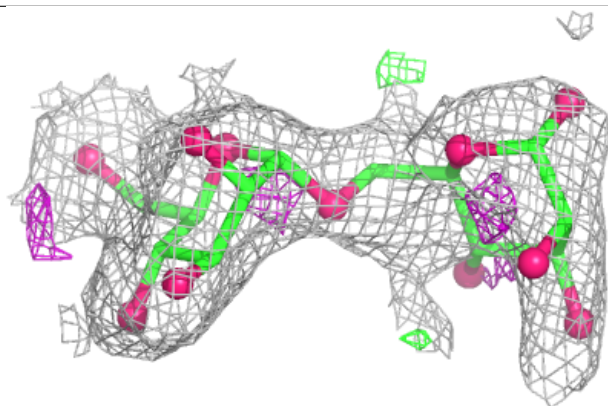
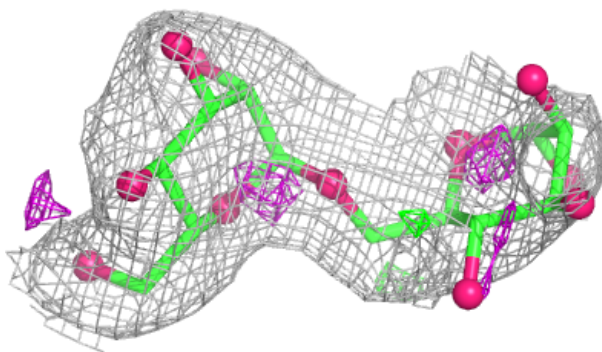


Electron density around Chain K:

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

**Electron density around Chain L:**

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

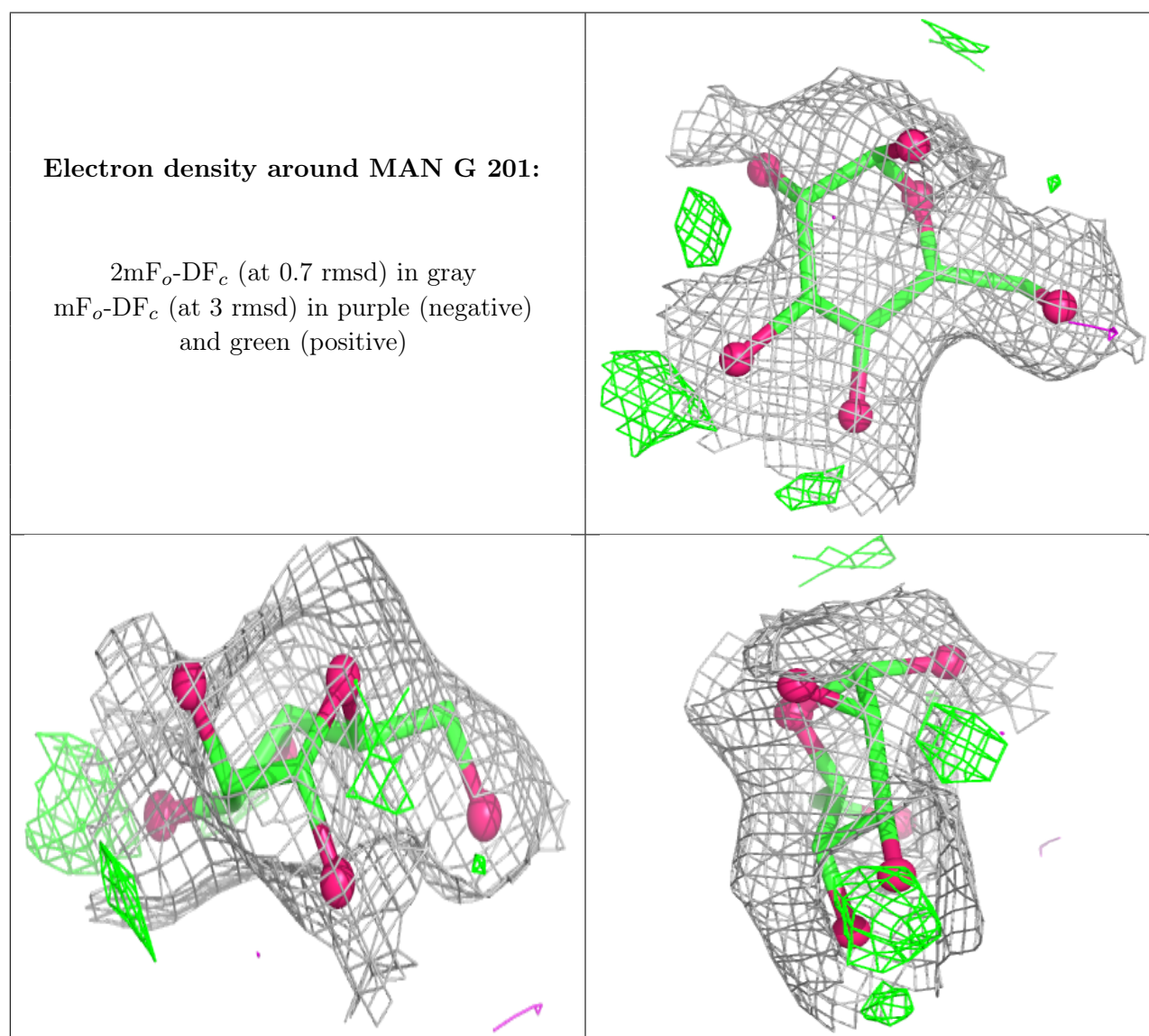


6.4 Ligands ⓘ

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.

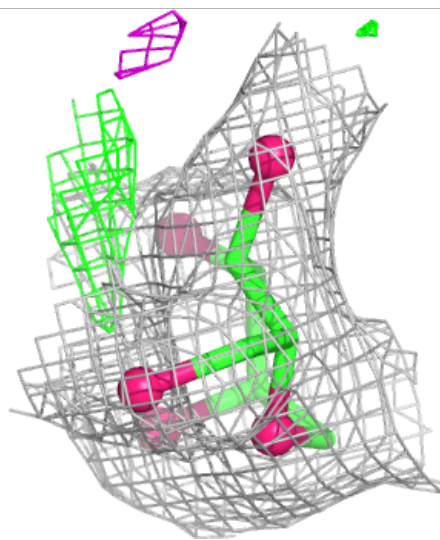
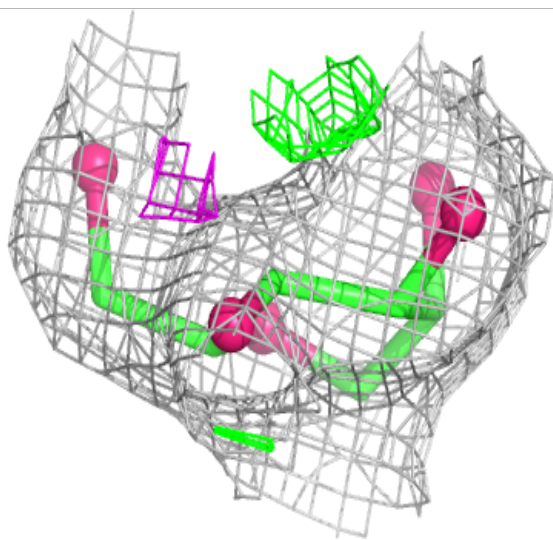
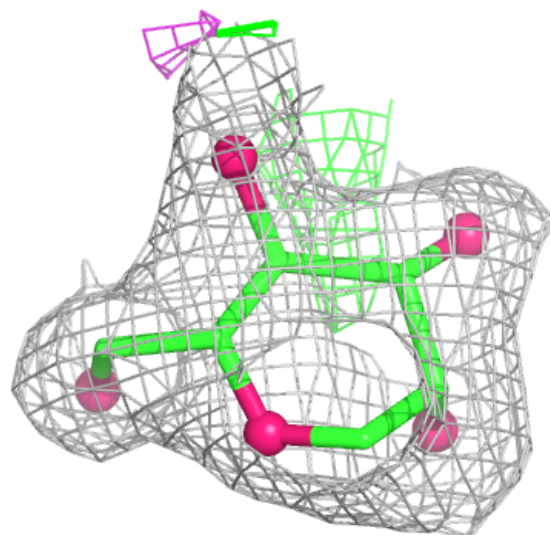
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
5	MAN	G	201	12/12	0.81	0.18	41,53,61,77	0
5	MAN	D	305	11/12	0.92	0.09	29,31,35,38	0
6	GOL	D	301	6/6	0.92	0.10	27,31,39,39	0

The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.



Electron density around MAN D 305:

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



6.5 Other polymers [i](#)

There are no such residues in this entry.