



Full wwPDB X-ray Structure Validation Report ⓘ

Aug 7, 2020 – 07:02 PM BST

PDB ID : 5AVA
Title : Crystal structure of PHA-E lectin in complex with bisected glycan
Authors : Nagae, M.; Yamaguchi, Y.
Deposited on : 2015-06-12
Resolution : 3.00 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.13.1
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.13.1

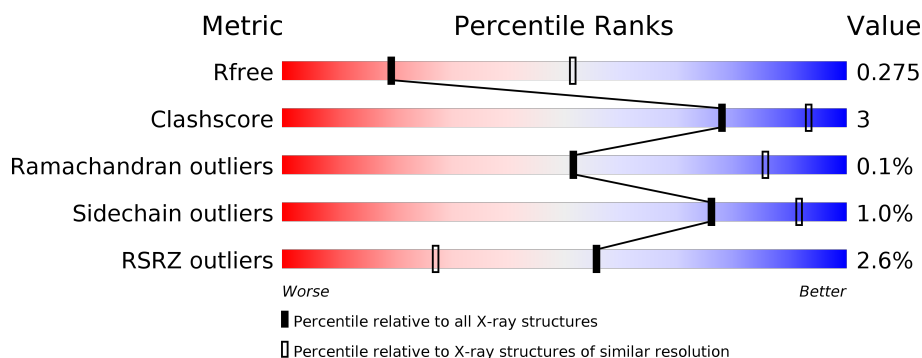
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.00 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	2092 (3.00-3.00)
Clashscore	141614	2416 (3.00-3.00)
Ramachandran outliers	138981	2333 (3.00-3.00)
Sidechain outliers	138945	2336 (3.00-3.00)
RSRZ outliers	127900	1990 (3.00-3.00)

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

Mol	Chain	Length	Quality of chain
1	A	275	<div> <div>2%</div> <div> <div></div> <div>78%</div> <div>8%</div> <div>13%</div> </div> </div>
1	B	275	<div> <div>2%</div> <div> <div></div> <div>78%</div> <div>8%</div> <div>13%</div> </div> </div>
1	C	275	<div> <div>2%</div> <div> <div></div> <div>81%</div> <div>5%</div> <div>13%</div> </div> </div>
1	D	275	<div> <div>2%</div> <div> <div></div> <div>80%</div> <div>6%</div> <div>13%</div> </div> </div>
1	E	275	<div> <div>%</div> <div> <div></div> <div>81%</div> <div>5%</div> <div>13%</div> </div> </div>
1	F	275	<div> <div>%</div> <div> <div></div> <div>77%</div> <div>9%</div> <div>13%</div> </div> </div>

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Mol	Chain	Length	Quality of chain
1	G	275	
1	H	275	
2	I	8	
2	J	8	
2	K	8	
2	L	8	
2	M	8	
2	N	8	
2	O	8	
2	P	8	

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
2	NAG	N	7	-	-	-	X

2 Entry composition [i](#)

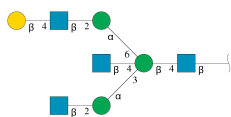
There are 5 unique types of molecules in this entry. The entry contains 15504 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Erythroagglutinin.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
1	A	238	Total	C	N	O	0	0	0
			1832	1169	300	363			
1	B	238	Total	C	N	O	0	0	0
			1832	1169	300	363			
1	C	238	Total	C	N	O	0	0	0
			1832	1169	300	363			
1	D	238	Total	C	N	O	0	0	0
			1832	1169	300	363			
1	E	238	Total	C	N	O	0	0	0
			1832	1169	300	363			
1	F	238	Total	C	N	O	0	0	0
			1832	1169	300	363			
1	G	238	Total	C	N	O	0	0	0
			1832	1169	300	363			
1	H	238	Total	C	N	O	0	0	0
			1832	1169	300	363			

- Molecule 2 is an oligosaccharide called beta-D-galactopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-2)-alpha-D-mannopyranose-(1-6)-[2-acetamido-2-deoxy-beta-D-glucopyranose-(1-2)-alpha-D-mannopyranose-(1-3)]|2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	I	8	Total	C	N	O	0	0	0
			100	56	4	40			
2	J	8	Total	C	N	O	0	0	0
			100	56	4	40			
2	K	8	Total	C	N	O	0	0	0
			100	56	4	40			

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	L	8	Total 100	C 56	N 4	O 40	0	0	0
2	M	8	Total 100	C 56	N 4	O 40	0	0	0
2	N	8	Total 100	C 56	N 4	O 40	0	0	0
2	O	8	Total 100	C 56	N 4	O 40	0	0	0
2	P	8	Total 100	C 56	N 4	O 40	0	0	0

- Molecule 3 is MANGANESE (II) ION (three-letter code: MN) (formula: Mn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	G	1	Total 1	Mn 1	0	0
3	D	1	Total 1	Mn 1	0	0
3	E	1	Total 1	Mn 1	0	0
3	H	1	Total 1	Mn 1	0	0
3	B	1	Total 1	Mn 1	0	0
3	C	1	Total 1	Mn 1	0	0
3	A	1	Total 1	Mn 1	0	0
3	F	1	Total 1	Mn 1	0	0

- Molecule 4 is CALCIUM ION (three-letter code: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	G	1	Total 1	Ca 1	0	0
4	D	1	Total 1	Ca 1	0	0
4	E	1	Total 1	Ca 1	0	0
4	H	1	Total 1	Ca 1	0	0

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	B	1	Total 1	Ca 1	0	0
4	C	1	Total 1	Ca 1	0	0
4	A	1	Total 1	Ca 1	0	0
4	F	1	Total 1	Ca 1	0	0

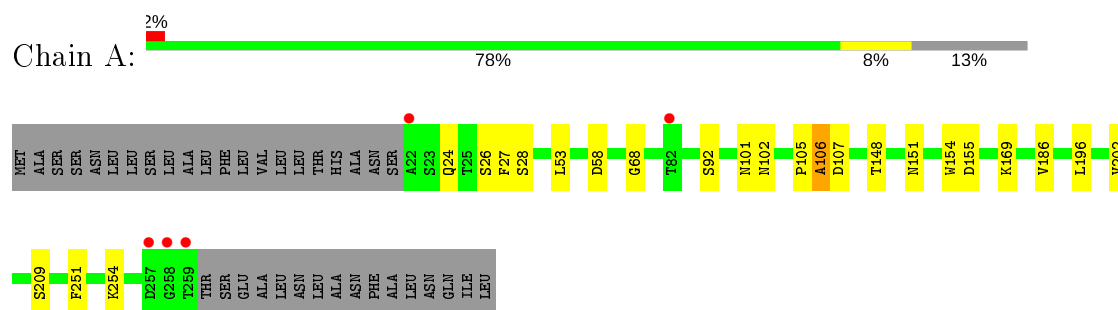
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	4	Total 4	O 4	0	0
5	B	4	Total 4	O 4	0	0
5	C	4	Total 4	O 4	0	0
5	D	4	Total 4	O 4	0	0
5	E	4	Total 4	O 4	0	0
5	F	4	Total 4	O 4	0	0
5	G	4	Total 4	O 4	0	0
5	H	4	Total 4	O 4	0	0

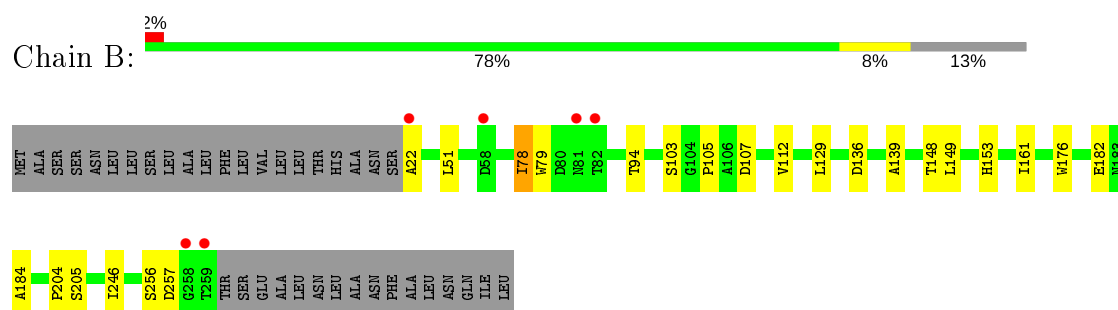
3 Residue-property plots [i](#)

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: Erythroagglutinin

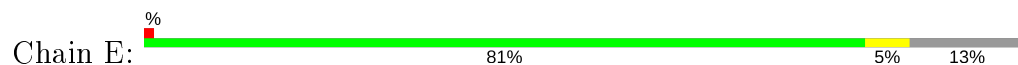


- Molecule 1: Erythroagglutinin



SER
GLU
ALA
LEU
LEU
LEU
LEU
ALA
ASN
PHE
ASN
ALA
ALA
LEU
LEU
LEU

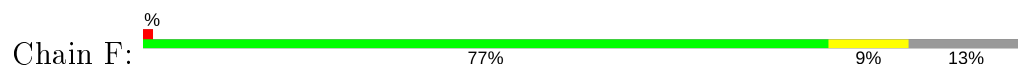
• Molecule 1: Erythroagglutinin



MET ALA SER SER ASN LEU LEU SER SER ALA LEU LEU PHE VAL LEU LEU THR THR HIS ALA ASN SER A22 L53 G68 Y72 S73 A74 I78 N102 A106 A143 L149 D164 S170 T173 E185 L196 V202 D257 G258 T259 THR SER GLU ALA

LEU
ASN
LEU
ALA
PHE
LEU
ALA
LEU
ASN
GLN
ILE
LEU

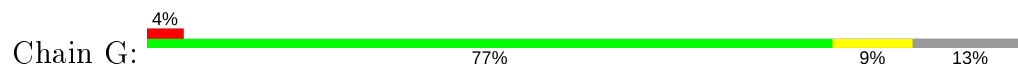
• Molecule 1: Erythroagglutinin



MET ALA SER SER ASN LEU LEU SER SER ALA LEU LEU PHE VAL LEU LEU THR THR HIS ALA ASN SER A22 S23 Q30 L51 R52 L53 G68 I78 W79 P105 A106 D107 V112 Q119 D122 L126 L129 A143 T148 L149 Y150 N151 D155 P156 K157

D164 L196 V226 T237 T246 S256 D257 G258 T259 THR SER SER GLU ALA LEU ASN SER A22 S23 PHE ALA LEU LEU ASN R52 L53 G68 I78 W79 P105 A106 D107 V112 Q119 D122 L126 L129 A143 T148 L149 Y150 N151 D155 P156 K157

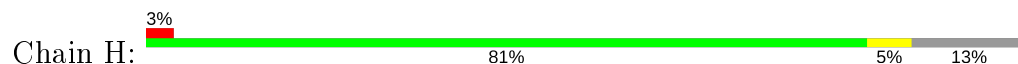
• Molecule 1: Erythroagglutinin



MET ALA SER SER ASN LEU LEU SER SER ALA LEU LEU PHE VAL LEU LEU THR THR HIS ALA ASN SER A22 S23 Q24 L51 R52 L53 N57 D58 N59 G60 T63 G68 Y72 S73 A74 I78 N102 P105 A106 D107 V112 L113 L126 L129 F130 T148 L149

H153 P204 V228 A232 I236 N240 N244 D245 K254 D257 G258 T259 THR SER GLU ALA LEU ASN SER A22 S23 PHE ALA LEU LEU ASN R52 L53 G68 I78 W79 P105 A106 D107 V112 Q119 D122 L126 L129 F130 T148 L149

• Molecule 1: Erythroagglutinin

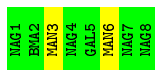


MET ALA SER SER ASN LEU LEU SER SER ALA LEU LEU PHE VAL LEU LEU THR THR HIS ALA ASN SER A22 L51 D58 Y72 S73 A74 A85 P105 V112 L126 L129 F130 T148 L149 H153 V202 Y203 P204 S209 G239 T243 T246 D257


G258 T259 THR SER GLU ALA LEU ASN SER A22 L51 D58 Y72 S73 A74 A85 P105 V112 L126 L129 F130 T148 L149 H153 V202 Y203 P204 S209 G239 T243 T246 D257

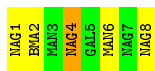
• Molecule 2: beta-D-galactopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-2)-alpha-D-mannopyranose-(1-6)-[2-acetamido-2-deoxy-beta-D-glucopyranose-(1-2)-alpha-D-mannopyranose-(1-3)]|2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain I:  75% 25%



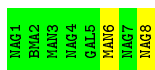
- Molecule 2: beta-D-galactopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-2)-alpha-D-mannopyranose-(1-6)-[2-acetamido-2-deoxy-beta-D-glucopyranose-(1-2)-alpha-D-mannopyranose-(1-3)][2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain J:  38% 50% 13%



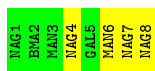
- Molecule 2: beta-D-galactopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-2)-alpha-D-mannopyranose-(1-6)-[2-acetamido-2-deoxy-beta-D-glucopyranose-(1-2)-alpha-D-mannopyranose-(1-3)][2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain K:  75% 25%




- Molecule 2: beta-D-galactopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-2)-alpha-D-mannopyranose-(1-6)-[2-acetamido-2-deoxy-beta-D-glucopyranose-(1-2)-alpha-D-mannopyranose-(1-3)][2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain L:  50% 50%



- Molecule 2: beta-D-galactopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-2)-alpha-D-mannopyranose-(1-6)-[2-acetamido-2-deoxy-beta-D-glucopyranose-(1-2)-alpha-D-mannopyranose-(1-3)][2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain M:  13% 88%




- Molecule 2: beta-D-galactopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-2)-alpha-D-mannopyranose-(1-6)-[2-acetamido-2-deoxy-beta-D-glucopyranose-(1-2)-alpha-D-mannopyranose-(1-3)][2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain N:  50% 50%



- Molecule 2: beta-D-galactopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-2)-alpha-D-mannopyranose-(1-6)-[2-acetamido-2-deoxy-beta-D-glucopyranose-(1-2)-alpha-D-mannopyranose-(1-3)][2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain O:  25% 38% 38%



- Molecule 2: beta-D-galactopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-2)-alpha-D-mannopyranose-(1-6)-[2-acetamido-2-deoxy-beta-D-glucopyranose-(1-2)-alpha-D-mannopyranose-(1-3)][2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain P:  50% 38% 13%



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	95.21Å 122.84Å 97.66Å 90.00° 90.66° 90.00°	Depositor
Resolution (Å)	100.00 – 3.00 31.73 – 3.00	Depositor EDS
% Data completeness (in resolution range)	99.8 (100.00-3.00) 99.9 (31.73-3.00)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.07	Depositor
$\langle I/\sigma(I) \rangle$ ¹	5.24 (at 3.00Å)	Xtriage
Refinement program	REFMAC 5.8.0049	Depositor
R, R_{free}	0.229 , 0.274 0.229 , 0.275	Depositor DCC
R_{free} test set	2268 reflections (5.03%)	wwPDB-VP
Wilson B-factor (Å ²)	44.5	Xtriage
Anisotropy	0.122	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 11.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.003 for l,k,-h 0.019 for h,-k,-l 0.014 for l,-k,h	Xtriage
F_o, F_c correlation	0.89	EDS
Total number of atoms	15504	wwPDB-VP
Average B, all atoms (Å ²)	42.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.11% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: BMA, NAG, CA, MN, GAL, MAN

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	0.32	0/1874	0.52	0/2559
1	B	0.34	0/1874	0.52	0/2559
1	C	0.32	0/1874	0.52	0/2559
1	D	0.33	0/1874	0.53	0/2559
1	E	0.33	0/1874	0.53	0/2559
1	F	0.33	0/1874	0.50	0/2559
1	G	0.33	0/1874	0.51	0/2559
1	H	0.33	0/1874	0.51	0/2559
All	All	0.33	0/14992	0.52	0/20472

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	1832	0	1796	15	0
1	B	1832	0	1796	13	0
1	C	1832	0	1796	9	0
1	D	1832	0	1796	9	0
1	E	1832	0	1796	9	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	F	1832	0	1796	16	0
1	G	1832	0	1796	14	0
1	H	1832	0	1796	8	0
2	I	100	0	85	0	0
2	J	100	0	85	1	0
2	K	100	0	85	0	0
2	L	100	0	85	1	0
2	M	100	0	85	0	0
2	N	100	0	85	0	0
2	O	100	0	85	3	0
2	P	100	0	85	1	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
3	C	1	0	0	0	0
3	D	1	0	0	0	0
3	E	1	0	0	0	0
3	F	1	0	0	0	0
3	G	1	0	0	0	0
3	H	1	0	0	0	0
4	A	1	0	0	0	0
4	B	1	0	0	0	0
4	C	1	0	0	0	0
4	D	1	0	0	0	0
4	E	1	0	0	0	0
4	F	1	0	0	0	0
4	G	1	0	0	0	0
4	H	1	0	0	0	0
5	A	4	0	0	0	0
5	B	4	0	0	0	0
5	C	4	0	0	0	0
5	D	4	0	0	0	0
5	E	4	0	0	0	0
5	F	4	0	0	0	0
5	G	4	0	0	0	0
5	H	4	0	0	0	0
All	All	15504	0	15048	87	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 3.

All (87) 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:202:VAL:HG12	1:A:209:SER:HB3	1.70	0.72
1:A:58:ASP:HB3	1:F:119:GLN:HG3	1.78	0.65
1:F:196:LEU:HD22	1:H:204:PRO:HA	1.83	0.61
1:B:78:ILE:HG22	1:B:79:TRP:HD1	1.67	0.59
1:F:22:ALA:HB3	1:F:256:SER:HA	1.85	0.58
1:E:53:LEU:O	1:E:68:GLY:HA3	2.04	0.57
1:G:24:GLN:HG2	1:G:254:LYS:HG3	1.86	0.57
1:D:106:ALA:HB2	1:D:149:LEU:HD13	1.87	0.56
1:B:182:GLU:OE1	1:B:205:SER:OG	2.22	0.56
1:A:24:GLN:HG2	1:A:254:LYS:HG3	1.85	0.56
1:B:112:VAL:HG11	1:B:129:LEU:HD21	1.87	0.56
1:A:53:LEU:O	1:A:68:GLY:HA3	2.07	0.55
1:C:24:GLN:HG2	1:C:254:LYS:HG3	1.90	0.54
1:H:126:LEU:HD22	1:H:130:PHE:HB2	1.91	0.51
1:G:112:VAL:HG11	1:G:129:LEU:HD21	1.91	0.51
2:O:6:MAN:H2	2:O:8:NAG:H61	1.91	0.51
1:A:101:ASN:O	1:A:102:ASN:HB2	2.10	0.51
1:E:72:TYR:CE2	1:E:74:ALA:HB3	2.46	0.51
1:D:53:LEU:O	1:D:68:GLY:HA3	2.11	0.49
1:B:51:LEU:HB3	1:B:246:ILE:HB	1.95	0.48
1:G:105:PRO:HB2	1:G:148:THR:HB	1.94	0.48
1:F:78:ILE:HG22	1:F:79:TRP:HD1	1.78	0.48
1:D:105:PRO:HB2	1:D:148:THR:HB	1.95	0.48
1:F:112:VAL:HG11	1:F:129:LEU:HD21	1.96	0.48
1:F:151:ASN:H	1:F:155:ASP:HB2	1.78	0.48
1:E:196:LEU:HD22	1:G:204:PRO:HA	1.95	0.47
1:C:185:GLU:HB2	1:C:202:VAL:HG23	1.95	0.47
1:C:202:VAL:HG12	1:C:209:SER:HB3	1.96	0.47
1:D:185:GLU:HB2	1:D:202:VAL:HG23	1.95	0.47
1:A:26:SER:HA	1:A:251:PHE:O	2.15	0.47
1:C:53:LEU:O	1:C:68:GLY:HA3	2.14	0.47
1:F:150:TYR:CD1	1:F:157:LYS:HA	2.49	0.47
1:F:78:ILE:HG12	1:F:226:VAL:HG22	1.97	0.47
1:G:126:LEU:HD22	1:G:130:PHE:HB2	1.96	0.47
1:F:105:PRO:HB2	1:F:148:THR:HB	1.98	0.46
1:F:53:LEU:O	1:F:68:GLY:HA3	2.15	0.46
1:G:51:LEU:HB3	1:G:246:ILE:HB	1.97	0.46
1:B:161:ILE:HG13	1:B:176:TRP:HB2	1.97	0.46
1:D:161:ILE:HG13	1:D:176:TRP:HB2	1.97	0.46
1:H:51:LEU:HB3	1:H:246:ILE:HB	1.97	0.46
1:D:112:VAL:HG11	1:D:129:LEU:HD21	1.97	0.45
1:E:170:SER:HB2	1:E:173:THR:HG23	1.98	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:106:ALA:HA	1:A:107:ASP:HA	1.78	0.45
1:E:22:ALA:HA	1:F:30:GLN:HG3	1.99	0.45
1:G:72:TYR:CE2	1:G:74:ALA:HB3	2.52	0.45
1:B:22:ALA:HB3	1:B:256:SER:HA	1.99	0.45
1:B:153:HIS:HD2	2:J:4:NAG:H81	1.82	0.44
1:C:51:LEU:HB3	1:C:246:ILE:HB	1.99	0.44
1:B:136:ASP:HB3	1:B:139:ALA:HB2	2.00	0.44
1:F:106:ALA:HA	1:F:107:ASP:HA	1.77	0.44
1:H:105:PRO:HB2	1:H:148:THR:HB	1.99	0.44
1:G:240:ASN:ND2	2:O:5:GAL:O2	2.49	0.44
1:G:78:ILE:HA	1:G:78:ILE:HD13	1.83	0.43
1:F:112:VAL:HG12	1:F:143:ALA:HB2	1.99	0.43
1:A:92:SER:HA	1:A:186:VAL:O	2.18	0.43
1:H:202:VAL:HG12	1:H:209:SER:HB3	2.01	0.43
1:C:106:ALA:HA	1:C:107:ASP:HA	1.79	0.43
1:D:149:LEU:HD21	2:L:4:NAG:H4	2.00	0.43
1:A:196:LEU:HD22	1:B:204:PRO:HA	1.99	0.43
1:B:22:ALA:HB1	1:B:257:ASP:CG	2.39	0.43
1:A:58:ASP:CB	1:F:119:GLN:HG3	2.46	0.43
1:C:64:LEU:HD23	1:C:241:VAL:HB	1.99	0.43
1:H:149:LEU:HD21	2:P:4:NAG:H4	2.00	0.42
1:G:106:ALA:HB2	1:G:149:LEU:HD13	2.01	0.42
1:G:53:LEU:O	1:G:68:GLY:HA3	2.19	0.42
1:A:151:ASN:H	1:A:155:ASP:HB2	1.85	0.42
1:G:232:ALA:HB1	1:G:244:ASN:ND2	2.35	0.42
1:D:91:THR:HG22	1:D:188:ILE:HB	2.02	0.42
1:A:154:TRP:HB2	1:A:169:LYS:HG3	2.02	0.42
1:B:94:THR:HA	1:B:184:ALA:O	2.20	0.42
1:E:185:GLU:HB2	1:E:202:VAL:HG23	2.02	0.42
1:A:27:PHE:CZ	1:A:251:PHE:HB3	2.54	0.41
1:B:105:PRO:HB2	1:B:148:THR:HB	2.02	0.41
1:B:22:ALA:CB	1:B:257:ASP:H	2.33	0.41
1:D:61:GLU:HA	1:D:62:PRO:HD2	1.91	0.41
1:H:112:VAL:HG11	1:H:129:LEU:HD21	2.03	0.41
1:C:24:GLN:HE21	1:C:254:LYS:HD2	1.86	0.41
1:A:105:PRO:HB2	1:A:148:THR:HB	2.03	0.41
1:A:28:SER:O	1:C:23:SER:HA	2.21	0.41
1:G:153:HIS:CD2	2:O:4:NAG:H81	2.56	0.41
1:E:106:ALA:HB2	1:E:149:LEU:HD13	2.03	0.40
1:E:78:ILE:HD13	1:E:78:ILE:HA	1.95	0.40
1:F:51:LEU:HB3	1:F:246:ILE:HB	2.04	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:113:LEU:HG	1:G:228:VAL:HG12	2.03	0.40
1:H:72:TYR:CE2	1:H:74:ALA:HB3	2.57	0.40
1:E:143:ALA:HB3	1:E:164:ASP:HB2	2.04	0.40
1:F:122:ASP:HB3	1:F:126:LEU:HD12	2.02	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	236/275 (86%)	227 (96%)	8 (3%)	1 (0%)	34	72
1	B	236/275 (86%)	225 (95%)	11 (5%)	0	100	100
1	C	236/275 (86%)	226 (96%)	10 (4%)	0	100	100
1	D	236/275 (86%)	225 (95%)	11 (5%)	0	100	100
1	E	236/275 (86%)	222 (94%)	14 (6%)	0	100	100
1	F	236/275 (86%)	223 (94%)	13 (6%)	0	100	100
1	G	236/275 (86%)	227 (96%)	9 (4%)	0	100	100
1	H	236/275 (86%)	225 (95%)	11 (5%)	0	100	100
All	All	1888/2200 (86%)	1800 (95%)	87 (5%)	1 (0%)	51	85

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	106	ALA

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	207/238 (87%)	207 (100%)	0	100	100
1	B	207/238 (87%)	203 (98%)	4 (2%)	57	84
1	C	207/238 (87%)	204 (99%)	3 (1%)	67	88
1	D	207/238 (87%)	203 (98%)	4 (2%)	57	84
1	E	207/238 (87%)	207 (100%)	0	100	100
1	F	207/238 (87%)	204 (99%)	3 (1%)	67	88
1	G	207/238 (87%)	206 (100%)	1 (0%)	88	96
1	H	207/238 (87%)	206 (100%)	1 (0%)	88	96
All	All	1656/1904 (87%)	1640 (99%)	16 (1%)	76	91

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

Mol	Chain	Res	Type
1	B	78	ILE
1	B	103	SER
1	B	107	ASP
1	B	149	LEU
1	C	78	ILE
1	C	103	SER
1	C	202	VAL
1	D	59	ASN
1	D	103	SER
1	D	149	LEU
1	D	202	VAL
1	F	23	SER
1	F	78	ILE
1	F	164	ASP
1	G	107	ASP
1	H	243	THR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (22) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	119	GLN
1	A	244	ASN
1	B	153	HIS
1	C	24	GLN
1	C	153	HIS
1	D	138	ASN
1	D	240	ASN
1	D	244	ASN
1	E	244	ASN
1	F	59	ASN
1	F	138	ASN
1	F	153	HIS
1	F	240	ASN
1	F	244	ASN
1	G	30	GLN
1	G	138	ASN
1	G	153	HIS
1	G	240	ASN
1	H	119	GLN
1	H	153	HIS
1	H	240	ASN
1	H	244	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

56 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	$\# Z > 2$
2	NAG	J	1	2	14,14,15	0.53	0	17,19,21	0.96	1 (5%)
2	GAL	K	5	2	11,11,12	0.60	0	15,15,17	0.78	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	MAN	N	6	2	11,11,12	0.55	0	15,15,17	0.88	1 (6%)
2	MAN	K	6	2	11,11,12	0.53	0	15,15,17	0.93	1 (6%)
2	NAG	L	4	2	14,14,15	0.51	0	17,19,21	0.86	0
2	MAN	J	3	2	11,11,12	0.62	0	15,15,17	0.84	0
2	NAG	K	8	2	14,14,15	0.41	0	17,19,21	1.26	1 (5%)
2	NAG	J	4	2	14,14,15	0.54	0	17,19,21	0.95	1 (5%)
2	NAG	L	1	2	14,14,15	0.51	0	17,19,21	0.89	0
2	NAG	O	1	2	14,14,15	0.54	0	17,19,21	1.19	2 (11%)
2	MAN	M	6	2	11,11,12	0.52	0	15,15,17	1.11	1 (6%)
2	NAG	N	8	2	14,14,15	0.39	0	17,19,21	1.08	1 (5%)
2	NAG	O	7	2	14,14,15	0.46	0	17,19,21	0.92	1 (5%)
2	MAN	J	6	2	11,11,12	0.52	0	15,15,17	1.68	1 (6%)
2	NAG	N	4	2	14,14,15	0.49	0	17,19,21	0.81	0
2	NAG	M	7	2	14,14,15	0.49	0	17,19,21	0.91	1 (5%)
2	MAN	N	3	2	11,11,12	0.61	0	15,15,17	0.67	0
2	NAG	N	1	2	14,14,15	0.49	0	17,19,21	0.99	1 (5%)
2	GAL	I	5	2	11,11,12	0.61	0	15,15,17	0.71	0
2	GAL	O	5	2	11,11,12	0.60	0	15,15,17	0.88	1 (6%)
2	NAG	J	8	2	14,14,15	0.49	0	17,19,21	1.32	2 (11%)
2	GAL	J	5	2	11,11,12	0.63	0	15,15,17	0.76	0
2	MAN	I	3	2	11,11,12	0.70	0	15,15,17	0.98	1 (6%)
2	MAN	K	3	2	11,11,12	0.63	0	15,15,17	0.66	0
2	NAG	M	4	2	14,14,15	0.46	0	17,19,21	0.83	0
2	MAN	M	3	2	11,11,12	0.61	0	15,15,17	1.02	1 (6%)
2	NAG	O	4	2	14,14,15	0.51	0	17,19,21	0.93	0
2	NAG	I	4	2	14,14,15	0.47	0	17,19,21	0.97	0
2	MAN	L	6	2	11,11,12	0.51	0	15,15,17	1.22	1 (6%)
2	MAN	O	6	2	11,11,12	0.53	0	15,15,17	1.17	2 (13%)
2	NAG	N	7	2	14,14,15	0.49	0	17,19,21	1.03	0
2	NAG	P	4	2	14,14,15	0.55	0	17,19,21	1.10	3 (17%)
2	NAG	O	8	2	14,14,15	0.38	0	17,19,21	1.31	1 (5%)
2	NAG	K	1	2	14,14,15	0.50	0	17,19,21	0.99	0
2	MAN	I	6	2	11,11,12	0.49	0	15,15,17	1.10	1 (6%)
2	NAG	L	7	2	14,14,15	0.52	0	17,19,21	0.85	1 (5%)
2	MAN	P	3	2	11,11,12	0.64	0	15,15,17	0.93	0
2	NAG	I	1	2	14,14,15	0.54	0	17,19,21	0.94	0
2	NAG	I	8	2	14,14,15	0.47	0	17,19,21	0.74	0
2	GAL	M	5	2	11,11,12	0.66	0	15,15,17	0.98	1 (6%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	NAG	L	8	2	14,14,15	0.48	0	17,19,21	0.96	1 (5%)
2	NAG	P	8	2	14,14,15	0.47	0	17,19,21	1.10	1 (5%)
2	NAG	P	7	2	14,14,15	0.52	0	17,19,21	0.62	0
2	GAL	N	5	2	11,11,12	0.61	0	15,15,17	0.77	0
2	MAN	O	3	2	11,11,12	0.67	0	15,15,17	0.70	0
2	NAG	P	1	2	14,14,15	0.55	0	17,19,21	0.90	1 (5%)
2	NAG	K	7	2	14,14,15	0.51	0	17,19,21	0.68	0
2	NAG	I	7	2	14,14,15	0.48	0	17,19,21	0.85	0
2	NAG	J	7	2	14,14,15	0.54	0	17,19,21	0.70	0
2	GAL	P	5	2	11,11,12	0.61	0	15,15,17	0.91	2 (13%)
2	MAN	L	3	2	11,11,12	0.64	0	15,15,17	0.64	0
2	MAN	P	6	2	11,11,12	0.58	0	15,15,17	0.79	0
2	NAG	M	1	2	14,14,15	0.53	0	17,19,21	0.94	1 (5%)
2	GAL	L	5	2	11,11,12	0.60	0	15,15,17	0.76	0
2	NAG	M	8	2	14,14,15	0.39	0	17,19,21	1.39	1 (5%)
2	NAG	K	4	2	14,14,15	0.52	0	17,19,21	0.85	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	NAG	J	1	2	-	0/6/23/26	0/1/1/1
2	GAL	K	5	2	-	2/2/19/22	0/1/1/1
2	MAN	N	6	2	-	0/2/19/22	0/1/1/1
2	MAN	K	6	2	-	2/2/19/22	0/1/1/1
2	NAG	L	4	2	-	0/6/23/26	0/1/1/1
2	MAN	J	3	2	-	0/2/19/22	0/1/1/1
2	NAG	K	8	2	-	0/6/23/26	0/1/1/1
2	NAG	J	4	2	-	2/6/23/26	0/1/1/1
2	NAG	L	1	2	-	0/6/23/26	0/1/1/1
2	NAG	O	1	2	-	1/6/23/26	0/1/1/1
2	MAN	M	6	2	-	2/2/19/22	0/1/1/1
2	NAG	N	8	2	-	0/6/23/26	0/1/1/1
2	NAG	O	7	2	-	2/6/23/26	0/1/1/1
2	MAN	J	6	2	-	2/2/19/22	0/1/1/1
2	NAG	N	4	2	-	0/6/23/26	0/1/1/1
2	NAG	M	7	2	-	2/6/23/26	0/1/1/1
2	MAN	N	3	2	-	0/2/19/22	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	NAG	N	1	2	-	2/6/23/26	0/1/1/1
2	GAL	I	5	2	-	2/2/19/22	0/1/1/1
2	GAL	O	5	2	-	2/2/19/22	0/1/1/1
2	NAG	J	8	2	-	0/6/23/26	0/1/1/1
2	GAL	J	5	2	-	2/2/19/22	0/1/1/1
2	MAN	I	3	2	-	1/2/19/22	0/1/1/1
2	MAN	K	3	2	-	0/2/19/22	0/1/1/1
2	NAG	M	4	2	-	0/6/23/26	0/1/1/1
2	MAN	M	3	2	-	2/2/19/22	0/1/1/1
2	NAG	O	4	2	-	0/6/23/26	0/1/1/1
2	NAG	I	4	2	-	0/6/23/26	0/1/1/1
2	MAN	L	6	2	-	0/2/19/22	0/1/1/1
2	MAN	O	6	2	-	1/2/19/22	0/1/1/1
2	NAG	N	7	2	-	2/6/23/26	0/1/1/1
2	NAG	P	4	2	-	0/6/23/26	0/1/1/1
2	NAG	O	8	2	-	2/6/23/26	0/1/1/1
2	NAG	K	1	2	-	2/6/23/26	0/1/1/1
2	MAN	I	6	2	-	1/2/19/22	0/1/1/1
2	NAG	L	7	2	-	2/6/23/26	0/1/1/1
2	MAN	P	3	2	-	2/2/19/22	0/1/1/1
2	NAG	I	1	2	-	0/6/23/26	0/1/1/1
2	NAG	I	8	2	-	0/6/23/26	0/1/1/1
2	GAL	M	5	2	-	2/2/19/22	0/1/1/1
2	NAG	L	8	2	-	0/6/23/26	0/1/1/1
2	NAG	P	8	2	-	0/6/23/26	0/1/1/1
2	NAG	P	7	2	-	0/6/23/26	0/1/1/1
2	GAL	N	5	2	-	2/2/19/22	0/1/1/1
2	MAN	O	3	2	-	0/2/19/22	0/1/1/1
2	NAG	P	1	2	-	0/6/23/26	0/1/1/1
2	NAG	K	7	2	-	2/6/23/26	0/1/1/1
2	NAG	I	7	2	-	2/6/23/26	0/1/1/1
2	NAG	J	7	2	-	2/6/23/26	0/1/1/1
2	GAL	P	5	2	-	2/2/19/22	0/1/1/1
2	MAN	L	3	2	-	0/2/19/22	0/1/1/1
2	MAN	P	6	2	-	1/2/19/22	0/1/1/1
2	NAG	M	1	2	-	0/6/23/26	0/1/1/1
2	GAL	L	5	2	-	2/2/19/22	0/1/1/1
2	NAG	M	8	2	-	0/6/23/26	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	NAG	K	4	2	-	0/6/23/26	0/1/1/1

There are no bond length outliers.

All (35) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	J	6	MAN	C1-O5-C5	5.47	119.60	112.19
2	M	8	NAG	C1-O5-C5	4.58	118.40	112.19
2	O	8	NAG	C1-O5-C5	4.52	118.31	112.19
2	K	8	NAG	C1-O5-C5	3.92	117.51	112.19
2	J	8	NAG	C1-O5-C5	3.42	116.83	112.19
2	L	6	MAN	C1-O5-C5	3.36	116.74	112.19
2	I	6	MAN	C1-O5-C5	3.35	116.73	112.19
2	N	8	NAG	C1-O5-C5	3.22	116.55	112.19
2	M	6	MAN	C1-O5-C5	3.17	116.48	112.19
2	P	8	NAG	C1-O5-C5	3.15	116.46	112.19
2	N	6	MAN	C1-O5-C5	2.81	116.00	112.19
2	K	6	MAN	C1-O5-C5	2.80	115.99	112.19
2	O	1	NAG	O5-C5-C6	2.71	111.46	107.20
2	O	7	NAG	C1-O5-C5	2.69	115.84	112.19
2	P	4	NAG	C2-N2-C7	2.59	126.60	122.90
2	M	3	MAN	C1-O5-C5	2.59	115.70	112.19
2	O	6	MAN	C1-O5-C5	2.59	115.70	112.19
2	M	5	GAL	O5-C1-C2	-2.56	106.81	110.77
2	N	1	NAG	C1-O5-C5	2.46	115.52	112.19
2	L	8	NAG	C1-O5-C5	2.41	115.46	112.19
2	J	1	NAG	O5-C5-C6	2.41	110.98	107.20
2	M	7	NAG	C1-O5-C5	2.37	115.41	112.19
2	J	8	NAG	C6-C5-C4	-2.29	107.65	113.00
2	M	1	NAG	O5-C5-C6	2.28	110.78	107.20
2	P	5	GAL	O5-C1-C2	-2.28	107.26	110.77
2	O	6	MAN	O5-C1-C2	-2.27	107.27	110.77
2	J	4	NAG	C1-O5-C5	2.17	115.13	112.19
2	O	1	NAG	C4-C3-C2	2.15	114.17	111.02
2	P	5	GAL	C1-O5-C5	2.11	115.05	112.19
2	P	4	NAG	O7-C7-N2	2.11	125.83	121.95
2	I	3	MAN	C1-C2-C3	2.09	112.24	109.67
2	P	4	NAG	C1-O5-C5	2.08	115.02	112.19
2	P	1	NAG	C4-C3-C2	2.08	114.07	111.02
2	O	5	GAL	O5-C5-C6	2.07	110.45	107.20
2	L	7	NAG	O5-C5-C6	2.04	110.41	107.20

There are no chirality outliers.

All (53) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	N	1	NAG	O5-C5-C6-O6
2	L	7	NAG	O5-C5-C6-O6
2	O	8	NAG	O5-C5-C6-O6
2	J	7	NAG	C4-C5-C6-O6
2	O	8	NAG	C4-C5-C6-O6
2	O	5	GAL	O5-C5-C6-O6
2	O	7	NAG	O5-C5-C6-O6
2	J	7	NAG	O5-C5-C6-O6
2	L	7	NAG	C4-C5-C6-O6
2	K	5	GAL	O5-C5-C6-O6
2	K	7	NAG	O5-C5-C6-O6
2	L	5	GAL	O5-C5-C6-O6
2	N	1	NAG	C4-C5-C6-O6
2	J	5	GAL	O5-C5-C6-O6
2	N	5	GAL	O5-C5-C6-O6
2	I	7	NAG	C4-C5-C6-O6
2	K	6	MAN	C4-C5-C6-O6
2	J	5	GAL	C4-C5-C6-O6
2	O	7	NAG	C4-C5-C6-O6
2	K	5	GAL	C4-C5-C6-O6
2	N	5	GAL	C4-C5-C6-O6
2	J	6	MAN	O5-C5-C6-O6
2	I	5	GAL	O5-C5-C6-O6
2	P	5	GAL	O5-C5-C6-O6
2	M	6	MAN	C4-C5-C6-O6
2	I	7	NAG	O5-C5-C6-O6
2	M	5	GAL	O5-C5-C6-O6
2	P	3	MAN	C4-C5-C6-O6
2	K	7	NAG	C4-C5-C6-O6
2	N	7	NAG	C4-C5-C6-O6
2	K	6	MAN	O5-C5-C6-O6
2	M	7	NAG	C4-C5-C6-O6
2	J	4	NAG	C4-C5-C6-O6
2	M	3	MAN	C4-C5-C6-O6
2	L	5	GAL	C4-C5-C6-O6
2	O	5	GAL	C4-C5-C6-O6
2	M	7	NAG	O5-C5-C6-O6
2	N	7	NAG	O5-C5-C6-O6
2	P	3	MAN	O5-C5-C6-O6
2	I	3	MAN	C4-C5-C6-O6

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Mol	Chain	Res	Type	Atoms
2	J	4	NAG	O5-C5-C6-O6
2	O	6	MAN	O5-C5-C6-O6
2	M	3	MAN	O5-C5-C6-O6
2	K	1	NAG	C4-C5-C6-O6
2	M	6	MAN	O5-C5-C6-O6
2	J	6	MAN	C4-C5-C6-O6
2	I	6	MAN	C4-C5-C6-O6
2	I	5	GAL	C4-C5-C6-O6
2	O	1	NAG	C4-C5-C6-O6
2	P	5	GAL	C4-C5-C6-O6
2	K	1	NAG	O5-C5-C6-O6
2	M	5	GAL	C4-C5-C6-O6
2	P	6	MAN	C4-C5-C6-O6

There are no ring outliers.

7 monomers are involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	L	4	NAG	1	0
2	J	4	NAG	1	0
2	O	5	GAL	1	0
2	O	4	NAG	1	0
2	O	6	MAN	1	0
2	P	4	NAG	1	0
2	O	8	NAG	1	0

5.5 Carbohydrates [i](#)

64 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	I	1	2	14,14,15	0.54	0	17,19,21	0.94	0
2	BMA	I	2	2	11,11,12	0.64	0	15,15,17	0.76	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	MAN	I	3	2	11,11,12	0.70	0	15,15,17	0.98	1 (6%)
2	NAG	I	4	2	14,14,15	0.47	0	17,19,21	0.97	0
2	GAL	I	5	2	11,11,12	0.61	0	15,15,17	0.71	0
2	MAN	I	6	2	11,11,12	0.49	0	15,15,17	1.10	1 (6%)
2	NAG	I	7	2	14,14,15	0.48	0	17,19,21	0.85	0
2	NAG	I	8	2	14,14,15	0.47	0	17,19,21	0.74	0
2	NAG	J	1	2	14,14,15	0.53	0	17,19,21	0.96	1 (5%)
2	BMA	J	2	2	11,11,12	0.58	0	15,15,17	0.85	1 (6%)
2	MAN	J	3	2	11,11,12	0.62	0	15,15,17	0.84	0
2	NAG	J	4	2	14,14,15	0.54	0	17,19,21	0.95	1 (5%)
2	GAL	J	5	2	11,11,12	0.63	0	15,15,17	0.76	0
2	MAN	J	6	2	11,11,12	0.52	0	15,15,17	1.68	1 (6%)
2	NAG	J	7	2	14,14,15	0.54	0	17,19,21	0.70	0
2	NAG	J	8	2	14,14,15	0.49	0	17,19,21	1.32	2 (11%)
2	NAG	K	1	2	14,14,15	0.50	0	17,19,21	0.99	0
2	BMA	K	2	2	11,11,12	0.60	0	15,15,17	0.74	0
2	MAN	K	3	2	11,11,12	0.63	0	15,15,17	0.66	0
2	NAG	K	4	2	14,14,15	0.52	0	17,19,21	0.85	0
2	GAL	K	5	2	11,11,12	0.60	0	15,15,17	0.78	0
2	MAN	K	6	2	11,11,12	0.53	0	15,15,17	0.93	1 (6%)
2	NAG	K	7	2	14,14,15	0.51	0	17,19,21	0.68	0
2	NAG	K	8	2	14,14,15	0.41	0	17,19,21	1.26	1 (5%)
2	NAG	L	1	2	14,14,15	0.51	0	17,19,21	0.89	0
2	BMA	L	2	2	11,11,12	0.54	0	15,15,17	0.63	0
2	MAN	L	3	2	11,11,12	0.64	0	15,15,17	0.64	0
2	NAG	L	4	2	14,14,15	0.51	0	17,19,21	0.86	0
2	GAL	L	5	2	11,11,12	0.60	0	15,15,17	0.76	0
2	MAN	L	6	2	11,11,12	0.51	0	15,15,17	1.22	1 (6%)
2	NAG	L	7	2	14,14,15	0.52	0	17,19,21	0.85	1 (5%)
2	NAG	L	8	2	14,14,15	0.48	0	17,19,21	0.96	1 (5%)
2	NAG	M	1	2	14,14,15	0.53	0	17,19,21	0.94	1 (5%)
2	BMA	M	2	2	11,11,12	0.55	0	15,15,17	0.84	1 (6%)
2	MAN	M	3	2	11,11,12	0.61	0	15,15,17	1.02	1 (6%)
2	NAG	M	4	2	14,14,15	0.46	0	17,19,21	0.83	0
2	GAL	M	5	2	11,11,12	0.66	0	15,15,17	0.98	1 (6%)
2	MAN	M	6	2	11,11,12	0.52	0	15,15,17	1.11	1 (6%)
2	NAG	M	7	2	14,14,15	0.49	0	17,19,21	0.91	1 (5%)
2	NAG	M	8	2	14,14,15	0.39	0	17,19,21	1.39	1 (5%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	NAG	N	1	2	14,14,15	0.49	0	17,19,21	0.99	1 (5%)
2	BMA	N	2	2	11,11,12	0.55	0	15,15,17	0.74	1 (6%)
2	MAN	N	3	2	11,11,12	0.61	0	15,15,17	0.67	0
2	NAG	N	4	2	14,14,15	0.49	0	17,19,21	0.81	0
2	GAL	N	5	2	11,11,12	0.61	0	15,15,17	0.77	0
2	MAN	N	6	2	11,11,12	0.55	0	15,15,17	0.88	1 (6%)
2	NAG	N	7	2	14,14,15	0.49	0	17,19,21	1.03	0
2	NAG	N	8	2	14,14,15	0.39	0	17,19,21	1.08	1 (5%)
2	NAG	O	1	2	14,14,15	0.54	0	17,19,21	1.19	2 (11%)
2	BMA	O	2	2	11,11,12	0.57	0	15,15,17	0.87	0
2	MAN	O	3	2	11,11,12	0.67	0	15,15,17	0.70	0
2	NAG	O	4	2	14,14,15	0.51	0	17,19,21	0.93	0
2	GAL	O	5	2	11,11,12	0.60	0	15,15,17	0.88	1 (6%)
2	MAN	O	6	2	11,11,12	0.53	0	15,15,17	1.17	2 (13%)
2	NAG	O	7	2	14,14,15	0.46	0	17,19,21	0.92	1 (5%)
2	NAG	O	8	2	14,14,15	0.38	0	17,19,21	1.31	1 (5%)
2	NAG	P	1	2	14,14,15	0.55	0	17,19,21	0.90	1 (5%)
2	BMA	P	2	2	11,11,12	0.52	0	15,15,17	0.54	0
2	MAN	P	3	2	11,11,12	0.64	0	15,15,17	0.93	0
2	NAG	P	4	2	14,14,15	0.55	0	17,19,21	1.10	3 (17%)
2	GAL	P	5	2	11,11,12	0.61	0	15,15,17	0.91	2 (13%)
2	MAN	P	6	2	11,11,12	0.58	0	15,15,17	0.79	0
2	NAG	P	7	2	14,14,15	0.52	0	17,19,21	0.62	0
2	NAG	P	8	2	14,14,15	0.47	0	17,19,21	1.10	1 (5%)

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	I	1	2	-	0/6/23/26	0/1/1/1
2	BMA	I	2	2	-	0/2/19/22	0/1/1/1
2	MAN	I	3	2	-	1/2/19/22	0/1/1/1
2	NAG	I	4	2	-	0/6/23/26	0/1/1/1
2	GAL	I	5	2	-	2/2/19/22	0/1/1/1
2	MAN	I	6	2	-	1/2/19/22	0/1/1/1
2	NAG	I	7	2	-	2/6/23/26	0/1/1/1
2	NAG	I	8	2	-	0/6/23/26	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	NAG	J	1	2	-	0/6/23/26	0/1/1/1
2	BMA	J	2	2	-	0/2/19/22	0/1/1/1
2	MAN	J	3	2	-	0/2/19/22	0/1/1/1
2	NAG	J	4	2	-	2/6/23/26	0/1/1/1
2	GAL	J	5	2	-	2/2/19/22	0/1/1/1
2	MAN	J	6	2	-	2/2/19/22	0/1/1/1
2	NAG	J	7	2	-	2/6/23/26	0/1/1/1
2	NAG	J	8	2	-	0/6/23/26	0/1/1/1
2	NAG	K	1	2	-	2/6/23/26	0/1/1/1
2	BMA	K	2	2	-	0/2/19/22	0/1/1/1
2	MAN	K	3	2	-	0/2/19/22	0/1/1/1
2	NAG	K	4	2	-	0/6/23/26	0/1/1/1
2	GAL	K	5	2	-	2/2/19/22	0/1/1/1
2	MAN	K	6	2	-	2/2/19/22	0/1/1/1
2	NAG	K	7	2	-	2/6/23/26	0/1/1/1
2	NAG	K	8	2	-	0/6/23/26	0/1/1/1
2	NAG	L	1	2	-	0/6/23/26	0/1/1/1
2	BMA	L	2	2	-	0/2/19/22	0/1/1/1
2	MAN	L	3	2	-	0/2/19/22	0/1/1/1
2	NAG	L	4	2	-	0/6/23/26	0/1/1/1
2	GAL	L	5	2	-	2/2/19/22	0/1/1/1
2	MAN	L	6	2	-	0/2/19/22	0/1/1/1
2	NAG	L	7	2	-	2/6/23/26	0/1/1/1
2	NAG	L	8	2	-	0/6/23/26	0/1/1/1
2	NAG	M	1	2	-	0/6/23/26	0/1/1/1
2	BMA	M	2	2	-	0/2/19/22	0/1/1/1
2	MAN	M	3	2	-	2/2/19/22	0/1/1/1
2	NAG	M	4	2	-	0/6/23/26	0/1/1/1
2	GAL	M	5	2	-	2/2/19/22	0/1/1/1
2	MAN	M	6	2	-	2/2/19/22	0/1/1/1
2	NAG	M	7	2	-	2/6/23/26	0/1/1/1
2	NAG	M	8	2	-	0/6/23/26	0/1/1/1
2	NAG	N	1	2	-	2/6/23/26	0/1/1/1
2	BMA	N	2	2	-	0/2/19/22	0/1/1/1
2	MAN	N	3	2	-	0/2/19/22	0/1/1/1
2	NAG	N	4	2	-	0/6/23/26	0/1/1/1
2	GAL	N	5	2	-	2/2/19/22	0/1/1/1
2	MAN	N	6	2	-	0/2/19/22	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	NAG	N	7	2	-	2/6/23/26	0/1/1/1
2	NAG	N	8	2	-	0/6/23/26	0/1/1/1
2	NAG	O	1	2	-	1/6/23/26	0/1/1/1
2	BMA	O	2	2	-	0/2/19/22	0/1/1/1
2	MAN	O	3	2	-	0/2/19/22	0/1/1/1
2	NAG	O	4	2	-	0/6/23/26	0/1/1/1
2	GAL	O	5	2	-	2/2/19/22	0/1/1/1
2	MAN	O	6	2	-	1/2/19/22	0/1/1/1
2	NAG	O	7	2	-	2/6/23/26	0/1/1/1
2	NAG	O	8	2	-	2/6/23/26	0/1/1/1
2	NAG	P	1	2	-	0/6/23/26	0/1/1/1
2	BMA	P	2	2	-	0/2/19/22	0/1/1/1
2	MAN	P	3	2	-	2/2/19/22	0/1/1/1
2	NAG	P	4	2	-	0/6/23/26	0/1/1/1
2	GAL	P	5	2	-	2/2/19/22	0/1/1/1
2	MAN	P	6	2	-	1/2/19/22	0/1/1/1
2	NAG	P	7	2	-	0/6/23/26	0/1/1/1
2	NAG	P	8	2	-	0/6/23/26	0/1/1/1

There are no bond length outliers.

All (38) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	J	6	MAN	C1-O5-C5	5.47	119.60	112.19
2	M	8	NAG	C1-O5-C5	4.58	118.40	112.19
2	O	8	NAG	C1-O5-C5	4.52	118.31	112.19
2	K	8	NAG	C1-O5-C5	3.92	117.51	112.19
2	J	8	NAG	C1-O5-C5	3.42	116.83	112.19
2	L	6	MAN	C1-O5-C5	3.36	116.74	112.19
2	I	6	MAN	C1-O5-C5	3.35	116.73	112.19
2	N	8	NAG	C1-O5-C5	3.22	116.55	112.19
2	M	6	MAN	C1-O5-C5	3.17	116.48	112.19
2	P	8	NAG	C1-O5-C5	3.15	116.46	112.19
2	N	6	MAN	C1-O5-C5	2.81	116.00	112.19
2	K	6	MAN	C1-O5-C5	2.80	115.99	112.19
2	O	1	NAG	O5-C5-C6	2.71	111.46	107.20
2	O	7	NAG	C1-O5-C5	2.69	115.84	112.19
2	P	4	NAG	C2-N2-C7	2.59	126.60	122.90
2	M	3	MAN	C1-O5-C5	2.59	115.70	112.19
2	O	6	MAN	C1-O5-C5	2.59	115.70	112.19

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	M	5	GAL	O5-C1-C2	-2.56	106.81	110.77
2	N	1	NAG	C1-O5-C5	2.46	115.52	112.19
2	L	8	NAG	C1-O5-C5	2.41	115.46	112.19
2	J	1	NAG	O5-C5-C6	2.41	110.98	107.20
2	M	7	NAG	C1-O5-C5	2.37	115.41	112.19
2	J	8	NAG	C6-C5-C4	-2.29	107.65	113.00
2	M	1	NAG	O5-C5-C6	2.28	110.78	107.20
2	P	5	GAL	O5-C1-C2	-2.28	107.26	110.77
2	O	6	MAN	O5-C1-C2	-2.27	107.27	110.77
2	J	4	NAG	C1-O5-C5	2.17	115.13	112.19
2	O	1	NAG	C4-C3-C2	2.15	114.17	111.02
2	M	2	BMA	C1-C2-C3	2.11	112.26	109.67
2	P	5	GAL	C1-O5-C5	2.11	115.05	112.19
2	P	4	NAG	O7-C7-N2	2.11	125.83	121.95
2	I	3	MAN	C1-C2-C3	2.09	112.24	109.67
2	P	4	NAG	C1-O5-C5	2.08	115.02	112.19
2	P	1	NAG	C4-C3-C2	2.08	114.07	111.02
2	O	5	GAL	O5-C5-C6	2.07	110.45	107.20
2	J	2	BMA	C1-O5-C5	2.05	114.97	112.19
2	L	7	NAG	O5-C5-C6	2.04	110.41	107.20
2	N	2	BMA	C1-C2-C3	2.00	112.12	109.67

There are no chirality outliers.

All (53) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	N	1	NAG	O5-C5-C6-O6
2	L	7	NAG	O5-C5-C6-O6
2	O	8	NAG	O5-C5-C6-O6
2	J	7	NAG	C4-C5-C6-O6
2	O	8	NAG	C4-C5-C6-O6
2	O	7	NAG	O5-C5-C6-O6
2	O	5	GAL	O5-C5-C6-O6
2	J	7	NAG	O5-C5-C6-O6
2	L	7	NAG	C4-C5-C6-O6
2	K	5	GAL	O5-C5-C6-O6
2	K	7	NAG	O5-C5-C6-O6
2	L	5	GAL	O5-C5-C6-O6
2	N	1	NAG	C4-C5-C6-O6
2	J	5	GAL	O5-C5-C6-O6
2	N	5	GAL	O5-C5-C6-O6
2	I	7	NAG	C4-C5-C6-O6

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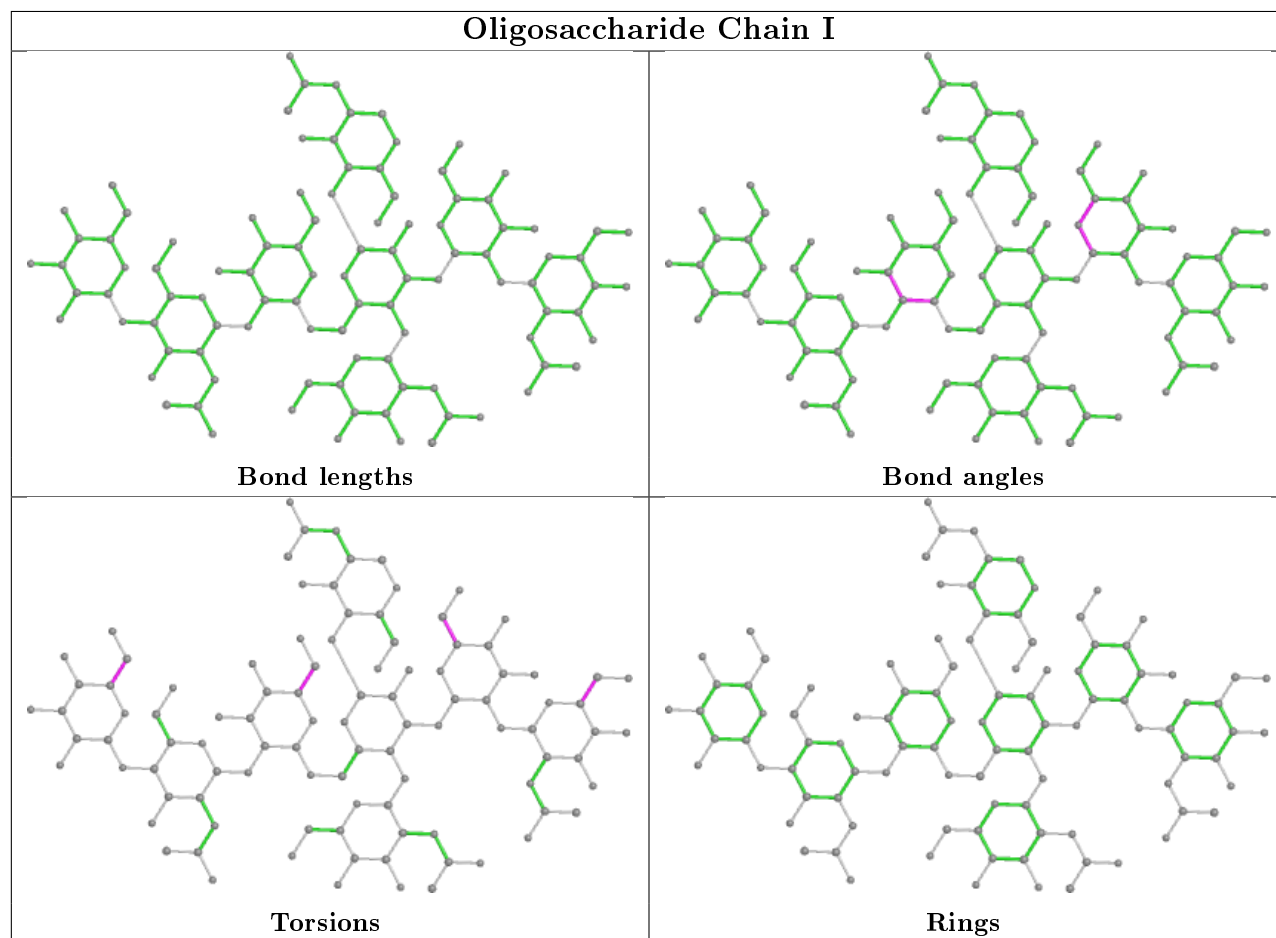
Mol	Chain	Res	Type	Atoms
2	K	6	MAN	C4-C5-C6-O6
2	O	7	NAG	C4-C5-C6-O6
2	J	5	GAL	C4-C5-C6-O6
2	K	5	GAL	C4-C5-C6-O6
2	N	5	GAL	C4-C5-C6-O6
2	P	5	GAL	O5-C5-C6-O6
2	J	6	MAN	O5-C5-C6-O6
2	I	5	GAL	O5-C5-C6-O6
2	M	6	MAN	C4-C5-C6-O6
2	I	7	NAG	O5-C5-C6-O6
2	M	5	GAL	O5-C5-C6-O6
2	P	3	MAN	C4-C5-C6-O6
2	K	7	NAG	C4-C5-C6-O6
2	N	7	NAG	C4-C5-C6-O6
2	K	6	MAN	O5-C5-C6-O6
2	M	7	NAG	C4-C5-C6-O6
2	J	4	NAG	C4-C5-C6-O6
2	M	3	MAN	C4-C5-C6-O6
2	L	5	GAL	C4-C5-C6-O6
2	O	5	GAL	C4-C5-C6-O6
2	M	7	NAG	O5-C5-C6-O6
2	N	7	NAG	O5-C5-C6-O6
2	P	3	MAN	O5-C5-C6-O6
2	I	3	MAN	C4-C5-C6-O6
2	J	4	NAG	O5-C5-C6-O6
2	O	6	MAN	O5-C5-C6-O6
2	M	3	MAN	O5-C5-C6-O6
2	K	1	NAG	C4-C5-C6-O6
2	M	6	MAN	O5-C5-C6-O6
2	J	6	MAN	C4-C5-C6-O6
2	I	6	MAN	C4-C5-C6-O6
2	I	5	GAL	C4-C5-C6-O6
2	O	1	NAG	C4-C5-C6-O6
2	P	5	GAL	C4-C5-C6-O6
2	K	1	NAG	O5-C5-C6-O6
2	M	5	GAL	C4-C5-C6-O6
2	P	6	MAN	C4-C5-C6-O6

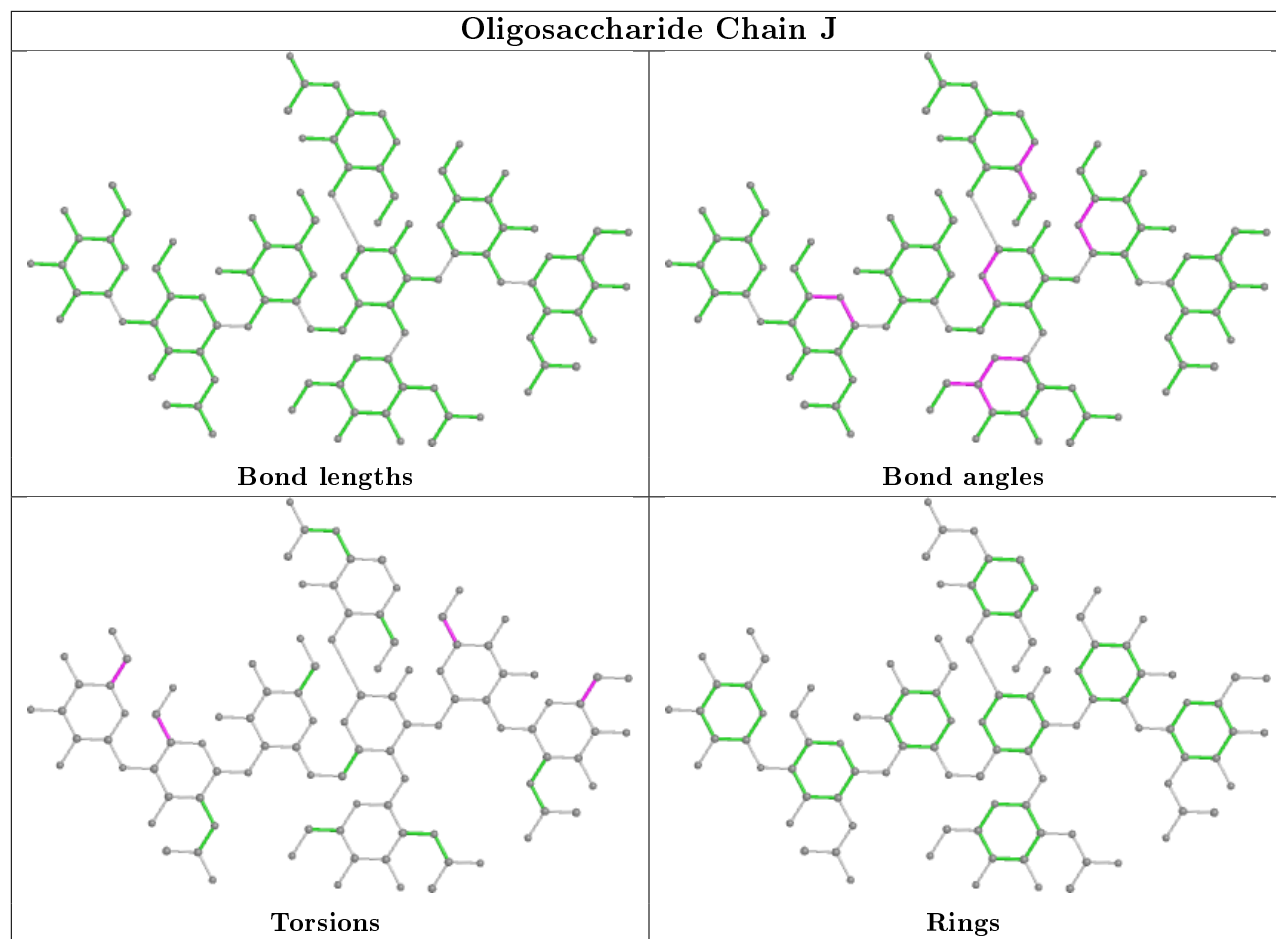
There are no ring outliers.

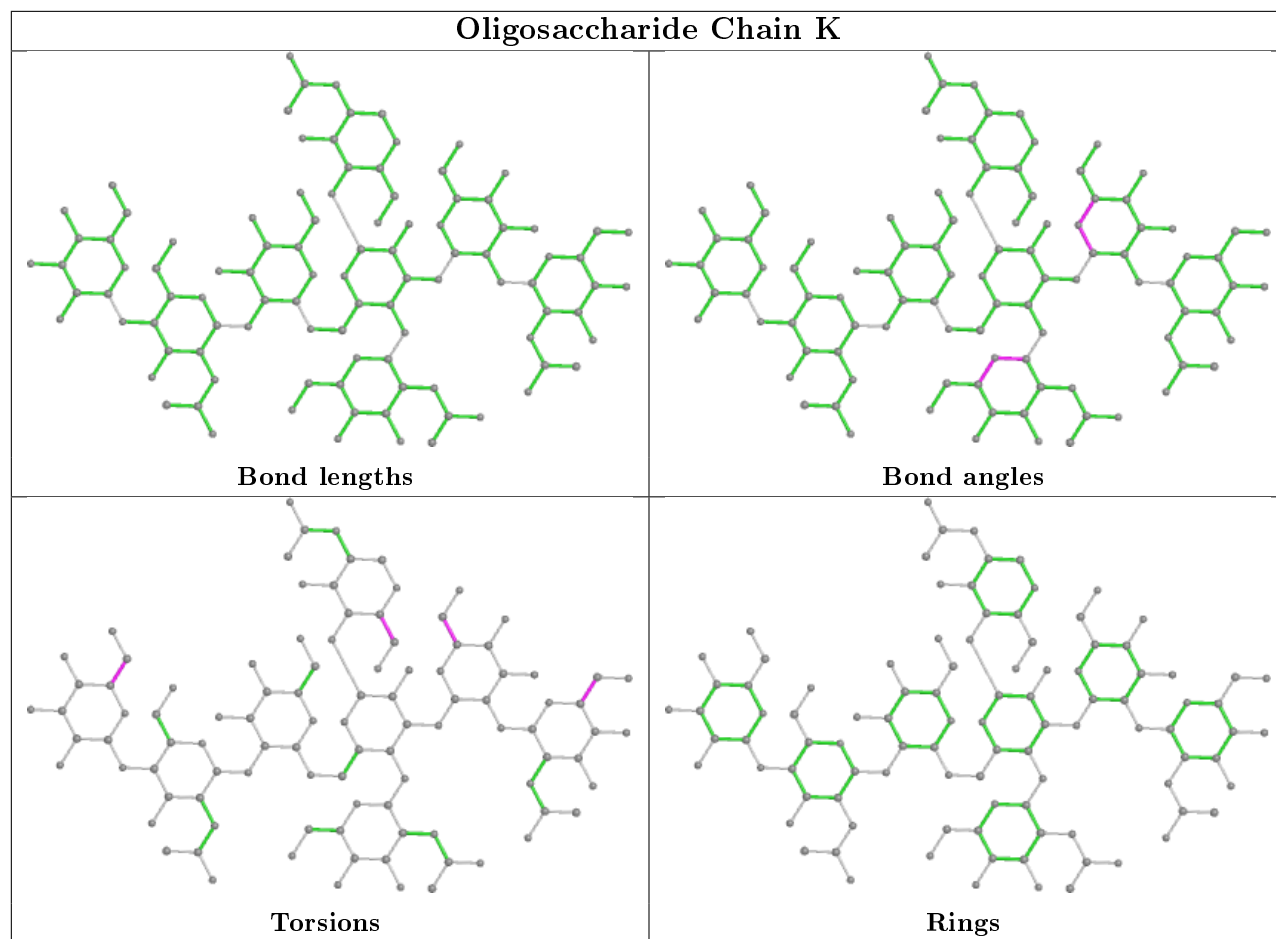
7 monomers are involved in 6 short contacts:

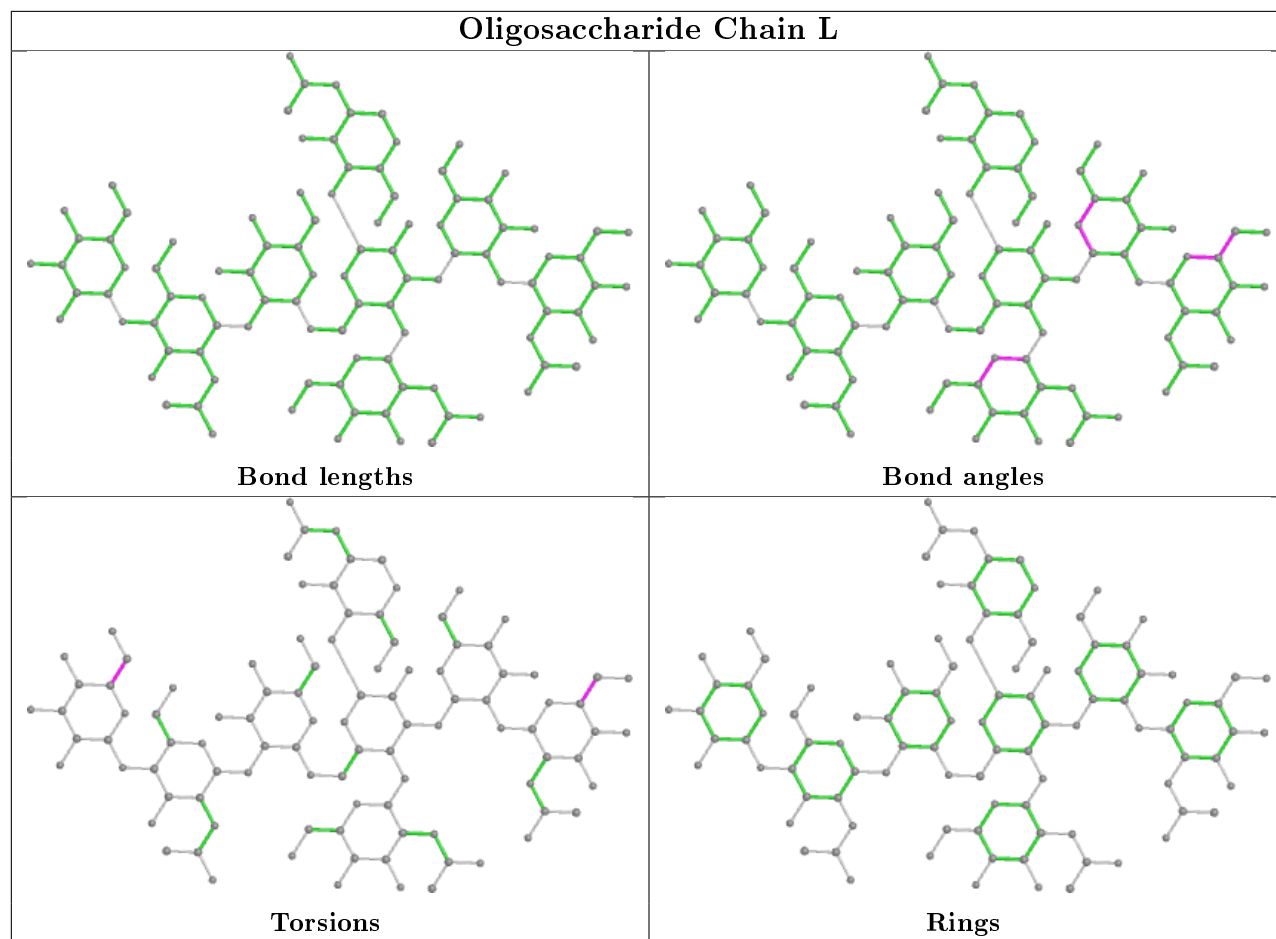
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	L	4	NAG	1	0
2	J	4	NAG	1	0
2	O	5	GAL	1	0
2	O	4	NAG	1	0
2	O	6	MAN	1	0
2	P	4	NAG	1	0
2	O	8	NAG	1	0

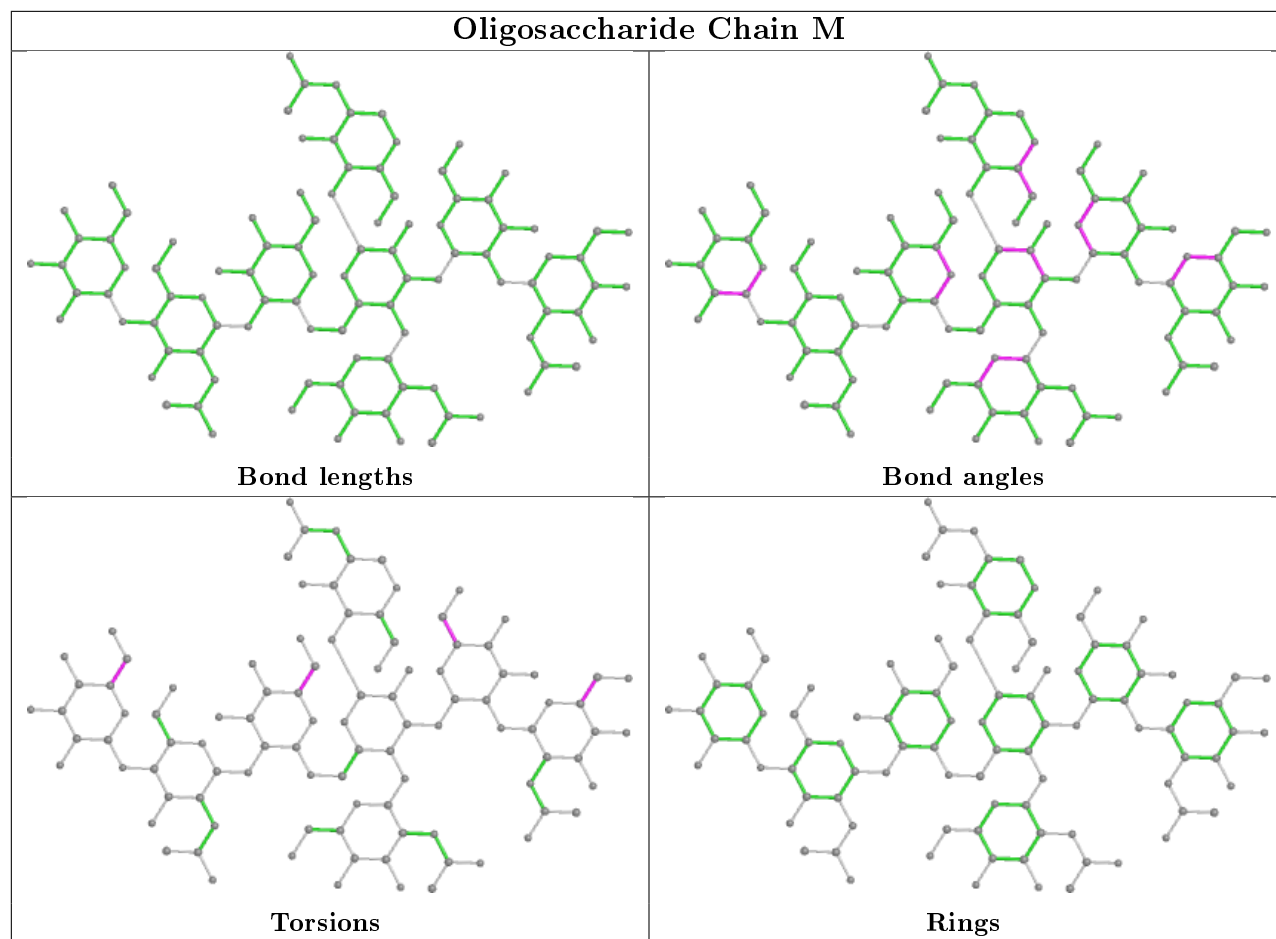
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for oligosaccharide.

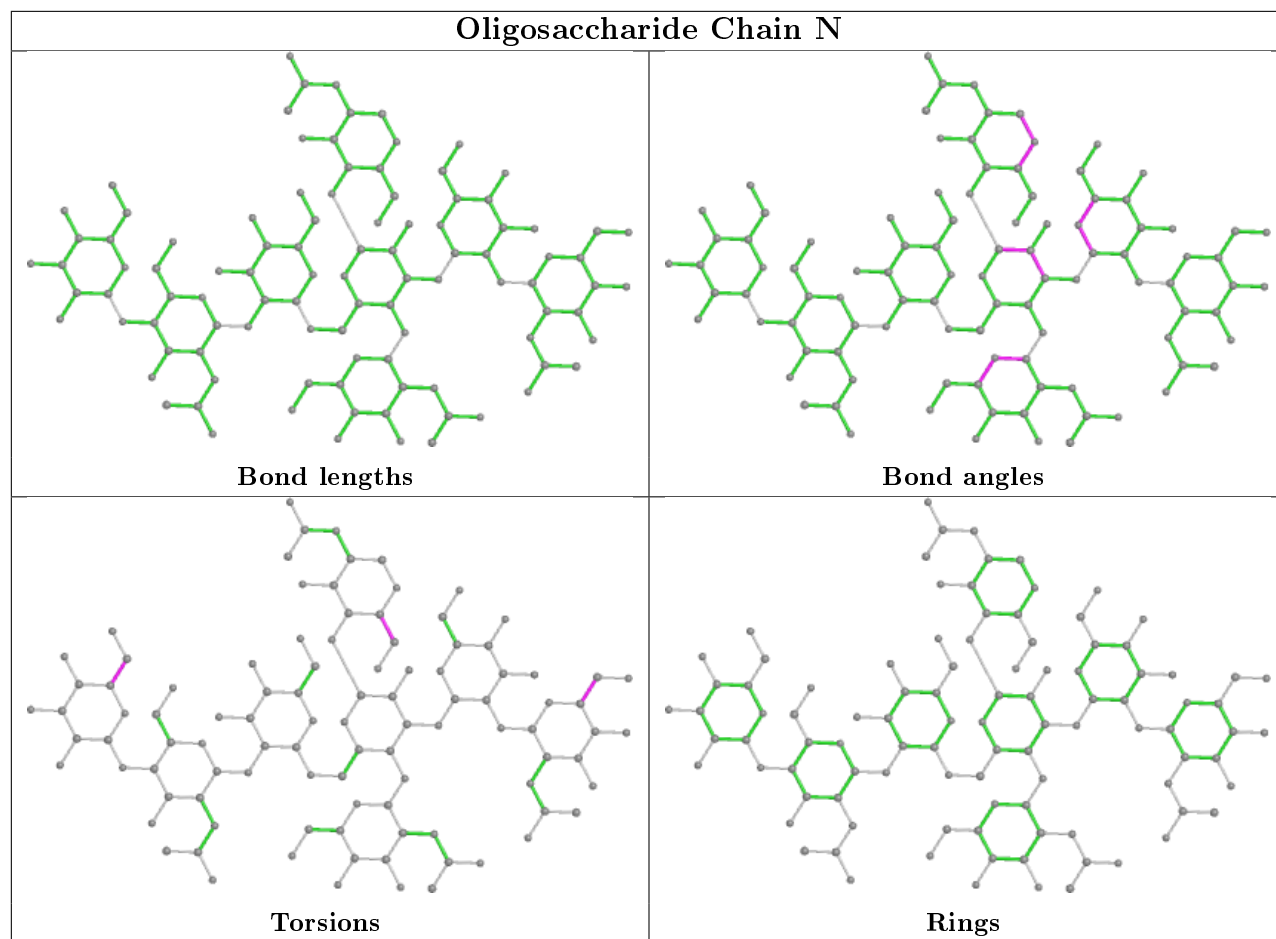


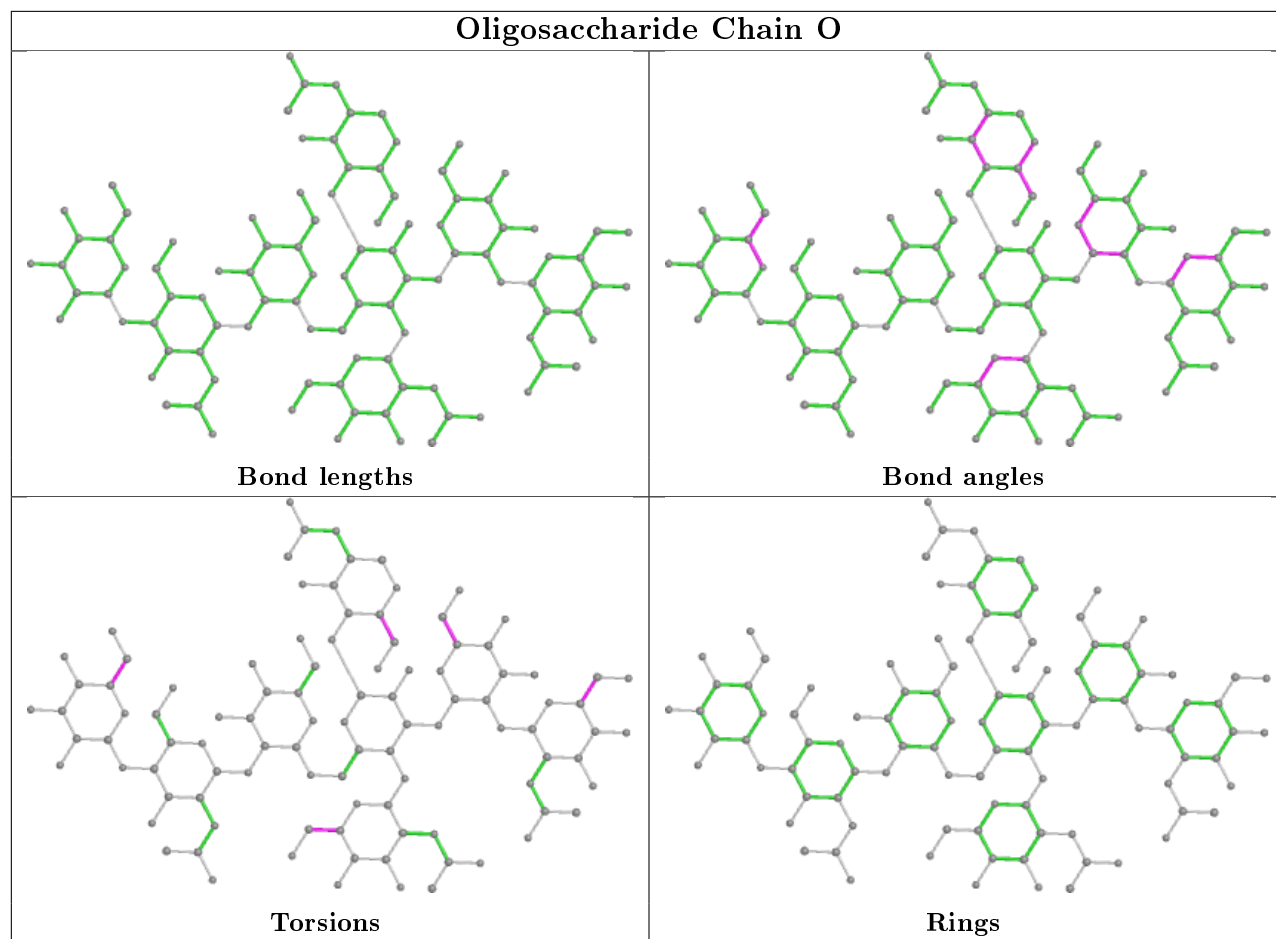


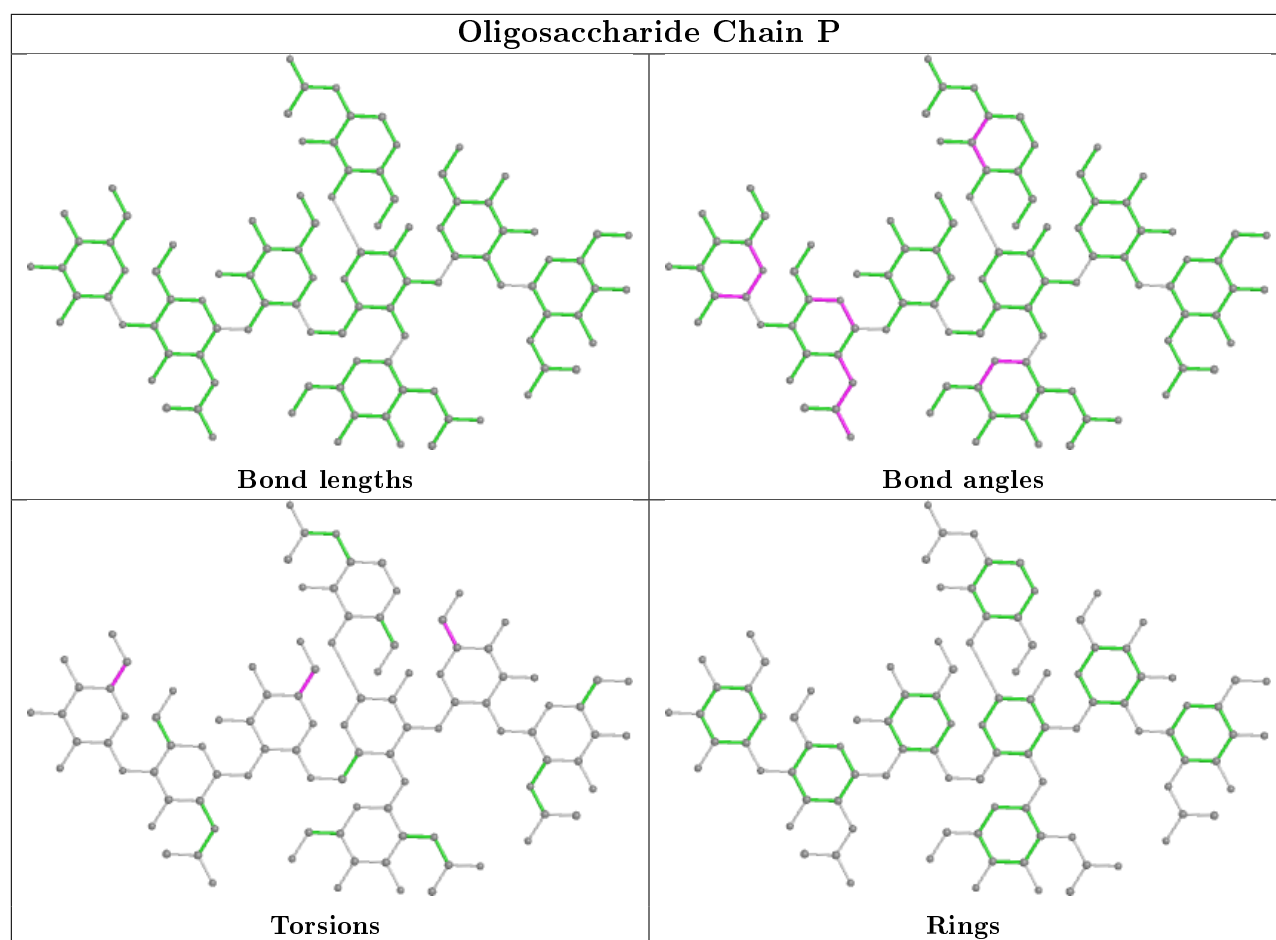












5.6 Ligand geometry [i](#)

Of 16 ligands modelled in this entry, 16 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues ⓘ

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	238/275 (86%)	-0.40	5 (2%)	63	34	20, 32, 50, 139	0
1	B	238/275 (86%)	-0.38	6 (2%)	57	29	22, 33, 51, 97	0
1	C	238/275 (86%)	-0.32	5 (2%)	63	34	23, 38, 68, 115	0
1	D	238/275 (86%)	-0.27	5 (2%)	63	34	24, 36, 61, 116	0
1	E	238/275 (86%)	-0.45	4 (1%)	70	41	19, 32, 51, 116	0
1	F	238/275 (86%)	-0.34	4 (1%)	70	41	20, 35, 62, 119	0
1	G	238/275 (86%)	0.06	12 (5%)	28	10	25, 50, 82, 131	0
1	H	238/275 (86%)	0.15	8 (3%)	45	19	30, 54, 83, 120	0
All	All	1904/2200 (86%)	-0.25	49 (2%)	56	27	19, 37, 73, 139	0

All (49) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	259	THR	10.8
1	D	259	THR	9.8
1	H	259	THR	9.6
1	E	259	THR	8.7
1	F	259	THR	8.3
1	C	259	THR	6.9
1	G	259	THR	6.0
1	E	258	GLY	5.8
1	A	258	GLY	4.6
1	G	59	ASN	4.4
1	G	258	GLY	4.3
1	B	259	THR	4.3
1	D	258	GLY	4.2
1	C	22	ALA	4.1
1	G	60	GLY	3.9
1	C	257	ASP	3.6

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Mol	Chain	Res	Type	RSRZ
1	H	258	GLY	3.6
1	G	22	ALA	3.5
1	H	257	ASP	3.4
1	C	258	GLY	3.2
1	A	22	ALA	3.1
1	F	258	GLY	3.1
1	B	58	ASP	3.0
1	H	58	ASP	2.9
1	F	22	ALA	2.8
1	D	257	ASP	2.8
1	B	258	GLY	2.8
1	B	22	ALA	2.7
1	A	257	ASP	2.6
1	G	257	ASP	2.6
1	G	153	HIS	2.6
1	D	22	ALA	2.6
1	E	102	ASN	2.5
1	H	22	ALA	2.4
1	G	236	ILE	2.4
1	B	82	THR	2.3
1	A	82	THR	2.3
1	H	153	HIS	2.2
1	D	59	ASN	2.2
1	H	85	ALA	2.2
1	H	239	GLY	2.2
1	G	58	ASP	2.2
1	B	81	ASN	2.2
1	G	63	THR	2.2
1	F	237	THR	2.1
1	C	83	THR	2.1
1	E	257	ASP	2.1
1	G	102	ASN	2.0
1	G	57	ASN	2.0

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

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(\AA^2)	Q<0.9
2	NAG	P	1	14/15	0.72	0.35	84,89,96,97	0
2	MAN	O	6	11/12	0.76	0.33	79,82,86,86	0
2	MAN	K	6	11/12	0.77	0.30	67,69,72,73	0
2	NAG	O	7	14/15	0.77	0.37	75,80,81,81	0
2	NAG	N	7	14/15	0.79	0.42	83,86,87,90	0
2	NAG	P	4	14/15	0.80	0.30	67,71,79,82	0
2	MAN	P	3	11/12	0.82	0.17	78,83,84,85	0
2	NAG	P	8	14/15	0.82	0.28	69,73,80,81	0
2	NAG	N	1	14/15	0.82	0.29	63,66,70,71	0
2	NAG	K	7	14/15	0.82	0.24	62,67,71,72	0
2	MAN	P	6	11/12	0.82	0.36	77,80,82,82	0
2	NAG	I	7	14/15	0.83	0.30	66,68,72,73	0
2	NAG	L	7	14/15	0.83	0.35	69,72,74,76	0
2	NAG	K	1	14/15	0.84	0.25	57,59,61,61	0
2	NAG	O	1	14/15	0.84	0.23	55,60,66,66	0
2	MAN	M	6	11/12	0.84	0.28	66,70,74,76	0
2	NAG	J	1	14/15	0.85	0.21	45,48,49,51	0
2	MAN	L	6	11/12	0.85	0.30	66,68,70,71	0
2	MAN	N	6	11/12	0.85	0.37	74,76,77,82	0
2	MAN	O	3	11/12	0.85	0.21	65,68,69,72	0
2	NAG	M	1	14/15	0.85	0.19	48,54,57,58	0
2	GAL	O	5	11/12	0.86	0.24	64,65,66,66	0
2	NAG	J	7	14/15	0.87	0.36	59,61,63,64	0
2	NAG	L	1	14/15	0.88	0.29	53,58,59,60	0
2	GAL	P	5	11/12	0.88	0.21	64,66,69,69	0
2	NAG	M	7	14/15	0.88	0.26	71,71,72,74	0
2	NAG	N	8	14/15	0.88	0.34	55,59,61,62	0
2	GAL	N	5	11/12	0.89	0.20	54,57,58,59	0
2	MAN	N	3	11/12	0.89	0.18	53,55,57,58	0
2	NAG	P	7	14/15	0.89	0.40	82,84,84,84	0
2	MAN	J	6	11/12	0.90	0.30	50,51,52,56	0
2	NAG	K	8	14/15	0.90	0.20	48,49,51,53	0
2	NAG	I	1	14/15	0.90	0.23	40,42,43,43	0
2	GAL	K	5	11/12	0.90	0.21	47,49,50,50	0
2	MAN	M	3	11/12	0.90	0.22	47,49,50,50	0
2	NAG	O	8	14/15	0.90	0.29	57,58,60,60	0
2	MAN	I	6	11/12	0.90	0.24	55,56,57,62	0
2	MAN	K	3	11/12	0.91	0.18	49,52,53,54	0
2	NAG	M	8	14/15	0.91	0.24	43,44,45,46	0
2	NAG	O	4	14/15	0.92	0.24	58,61,62,64	0
2	GAL	J	5	11/12	0.92	0.23	40,42,43,44	0
2	NAG	L	8	14/15	0.92	0.26	51,53,61,63	0
2	NAG	N	4	14/15	0.92	0.18	55,58,62,63	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
2	NAG	J	8	14/15	0.92	0.29	35,38,40,40	0
2	NAG	I	8	14/15	0.93	0.23	39,40,43,45	0
2	GAL	M	5	11/12	0.93	0.20	40,41,41,41	0
2	MAN	I	3	11/12	0.93	0.21	37,38,38,39	0
2	GAL	I	5	11/12	0.93	0.17	31,32,33,34	0
2	NAG	L	4	14/15	0.93	0.18	45,47,51,51	0
2	NAG	K	4	14/15	0.93	0.20	45,47,48,49	0
2	MAN	J	3	11/12	0.94	0.19	35,37,38,38	0
2	MAN	L	3	11/12	0.94	0.14	48,50,52,53	0
2	NAG	M	4	14/15	0.95	0.17	42,44,45,46	0
2	NAG	I	4	14/15	0.96	0.21	34,36,40,41	0
2	GAL	L	5	11/12	0.96	0.21	40,41,43,43	0
2	NAG	J	4	14/15	0.97	0.20	33,34,35,38	0

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(\AA^2)	Q<0.9
2	NAG	P	1	14/15	0.72	0.35	84,89,96,97	0
2	MAN	O	6	11/12	0.76	0.33	79,82,86,86	0
2	MAN	K	6	11/12	0.77	0.30	67,69,72,73	0
2	NAG	O	7	14/15	0.77	0.37	75,80,81,81	0
2	NAG	N	7	14/15	0.79	0.42	83,86,87,90	0
2	NAG	P	4	14/15	0.80	0.30	67,71,79,82	0
2	MAN	P	3	11/12	0.82	0.17	78,83,84,85	0
2	NAG	P	8	14/15	0.82	0.28	69,73,80,81	0
2	NAG	N	1	14/15	0.82	0.29	63,66,70,71	0
2	NAG	K	7	14/15	0.82	0.24	62,67,71,72	0
2	MAN	P	6	11/12	0.82	0.36	77,80,82,82	0
2	NAG	I	7	14/15	0.83	0.30	66,68,72,73	0
2	NAG	L	7	14/15	0.83	0.35	69,72,74,76	0
2	MAN	M	6	11/12	0.84	0.28	66,70,74,76	0
2	NAG	O	1	14/15	0.84	0.23	55,60,66,66	0
2	NAG	K	1	14/15	0.84	0.25	57,59,61,61	0
2	BMA	P	2	11/12	0.85	0.17	76,79,83,84	0
2	MAN	O	3	11/12	0.85	0.21	65,68,69,72	0
2	NAG	J	1	14/15	0.85	0.21	45,48,49,51	0
2	MAN	N	6	11/12	0.85	0.37	74,76,77,82	0

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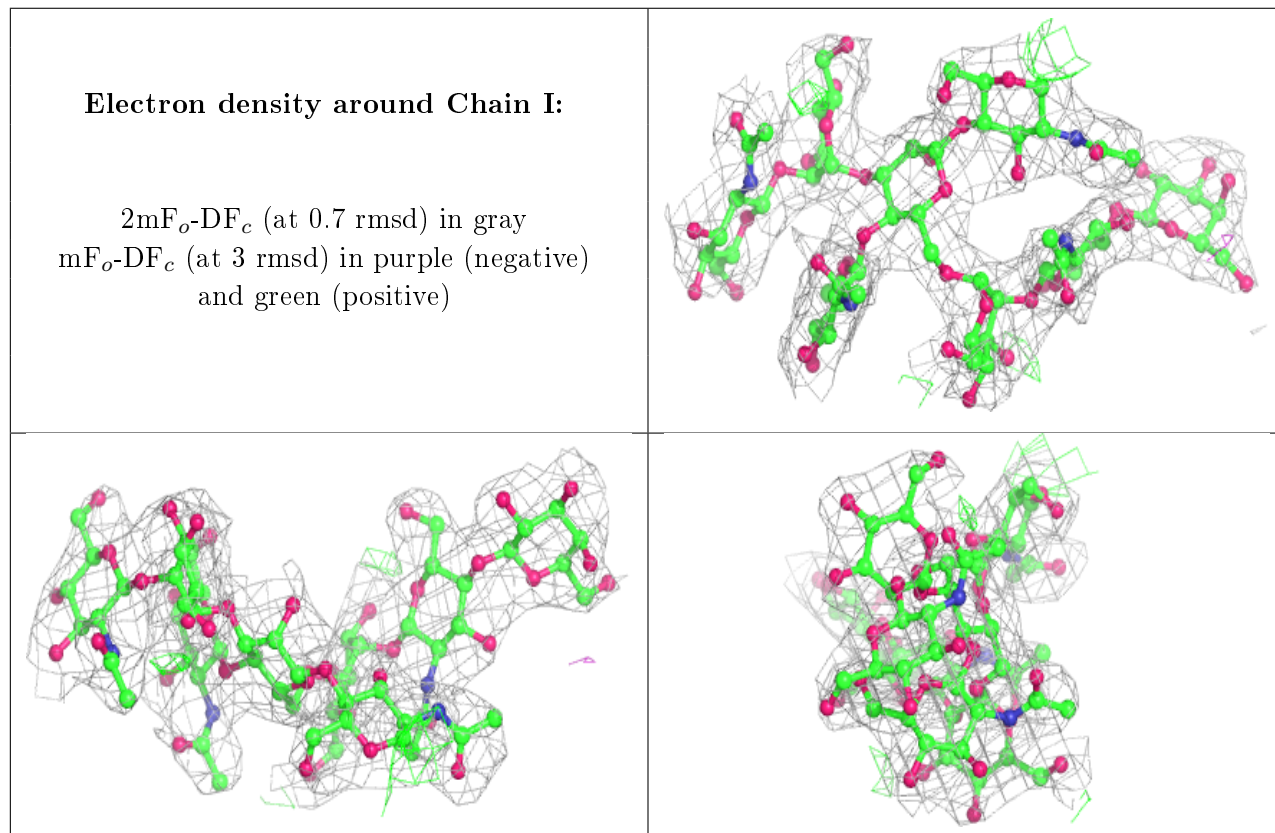
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
2	MAN	L	6	11/12	0.85	0.30	66,68,70,71	0
2	NAG	M	1	14/15	0.85	0.19	48,54,57,58	0
2	GAL	O	5	11/12	0.86	0.24	64,65,66,66	0
2	NAG	J	7	14/15	0.87	0.36	59,61,63,64	0
2	NAG	M	7	14/15	0.88	0.26	71,71,72,74	0
2	GAL	P	5	11/12	0.88	0.21	64,66,69,69	0
2	NAG	N	8	14/15	0.88	0.34	55,59,61,62	0
2	NAG	L	1	14/15	0.88	0.29	53,58,59,60	0
2	GAL	N	5	11/12	0.89	0.20	54,57,58,59	0
2	MAN	N	3	11/12	0.89	0.18	53,55,57,58	0
2	BMA	L	2	11/12	0.89	0.20	52,57,59,63	0
2	NAG	P	7	14/15	0.89	0.40	82,84,84,84	0
2	NAG	I	1	14/15	0.90	0.23	40,42,43,43	0
2	MAN	I	6	11/12	0.90	0.24	55,56,57,62	0
2	BMA	N	2	11/12	0.90	0.21	58,63,66,69	0
2	NAG	O	8	14/15	0.90	0.29	57,58,60,60	0
2	NAG	K	8	14/15	0.90	0.20	48,49,51,53	0
2	BMA	O	2	11/12	0.90	0.19	62,64,69,73	0
2	GAL	K	5	11/12	0.90	0.21	47,49,50,50	0
2	MAN	M	3	11/12	0.90	0.22	47,49,50,50	0
2	MAN	J	6	11/12	0.90	0.30	50,51,52,56	0
2	NAG	M	8	14/15	0.91	0.24	43,44,45,46	0
2	BMA	J	2	11/12	0.91	0.18	39,43,44,47	0
2	MAN	K	3	11/12	0.91	0.18	49,52,53,54	0
2	GAL	J	5	11/12	0.92	0.23	40,42,43,44	0
2	NAG	N	4	14/15	0.92	0.18	55,58,62,63	0
2	BMA	M	2	11/12	0.92	0.18	50,53,55,60	0
2	NAG	J	8	14/15	0.92	0.29	35,38,40,40	0
2	NAG	L	8	14/15	0.92	0.26	51,53,61,63	0
2	NAG	O	4	14/15	0.92	0.24	58,61,62,64	0
2	GAL	I	5	11/12	0.93	0.17	31,32,33,34	0
2	NAG	I	8	14/15	0.93	0.23	39,40,43,45	0
2	BMA	K	2	11/12	0.93	0.14	53,55,58,63	0
2	GAL	M	5	11/12	0.93	0.20	40,41,41,41	0
2	MAN	I	3	11/12	0.93	0.21	37,38,38,39	0
2	NAG	L	4	14/15	0.93	0.18	45,47,51,51	0
2	NAG	K	4	14/15	0.93	0.20	45,47,48,49	0
2	MAN	J	3	11/12	0.94	0.19	35,37,38,38	0
2	MAN	L	3	11/12	0.94	0.14	48,50,52,53	0
2	BMA	I	2	11/12	0.95	0.18	39,42,45,50	0
2	NAG	M	4	14/15	0.95	0.17	42,44,45,46	0
2	NAG	I	4	14/15	0.96	0.21	34,36,40,41	0

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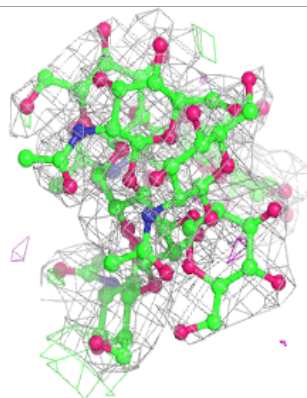
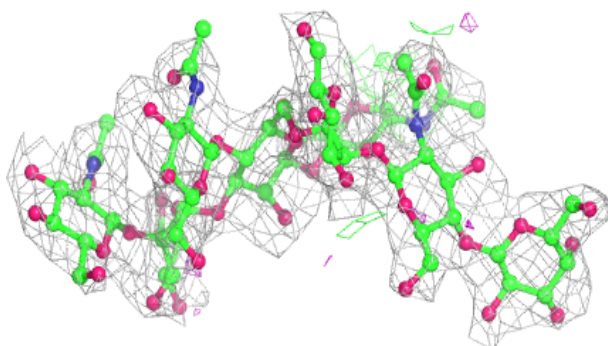
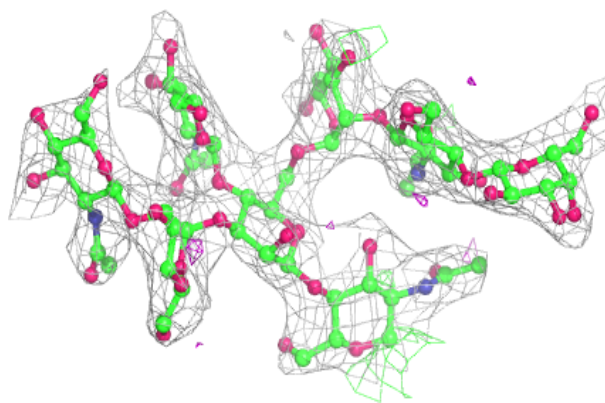
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
2	GAL	L	5	11/12	0.96	0.21	40,41,43,43	0
2	NAG	J	4	14/15	0.97	0.20	33,34,35,38	0

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.

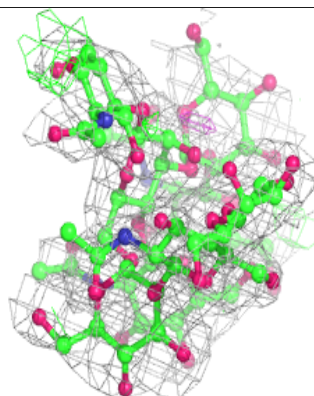
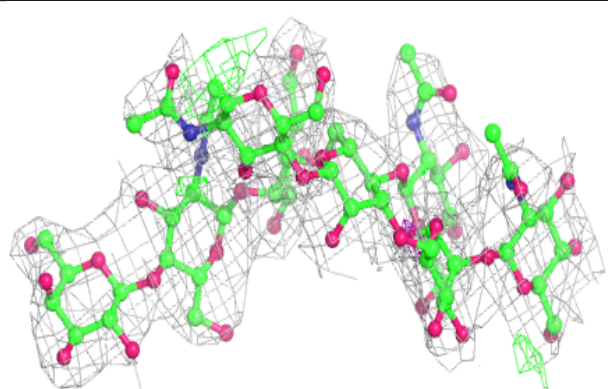
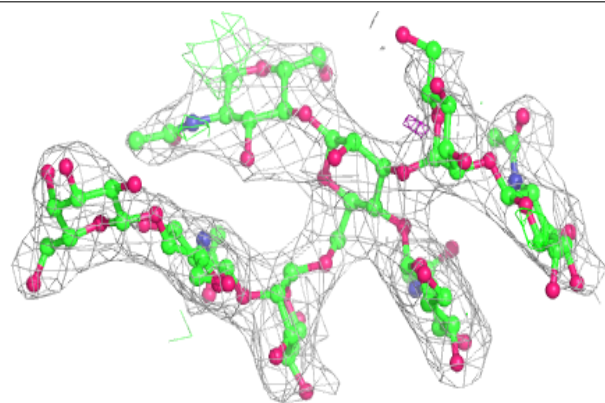


Electron density around Chain J:

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

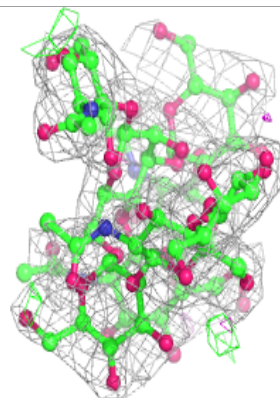
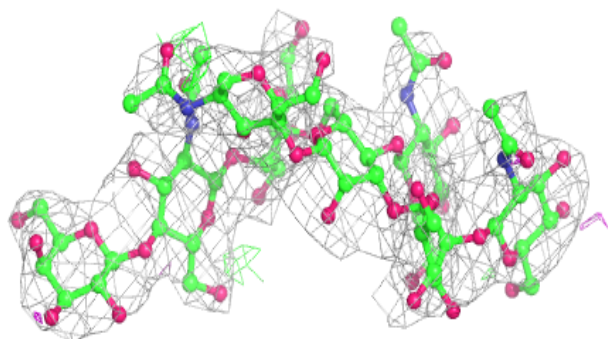
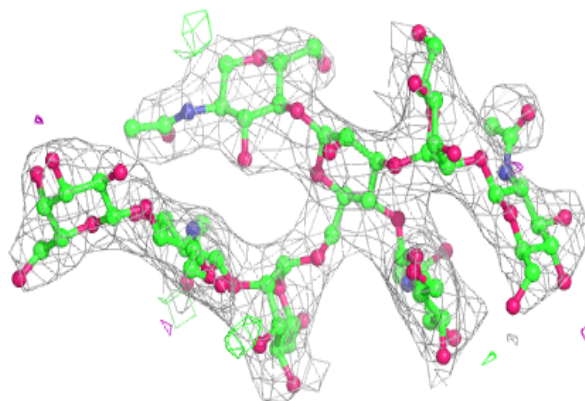
**Electron density around Chain K:**

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

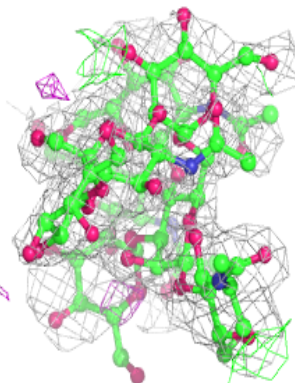
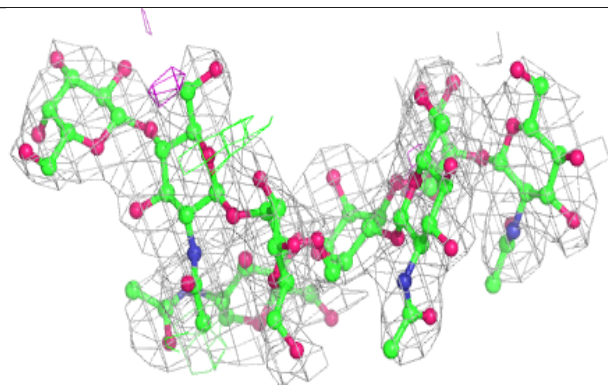
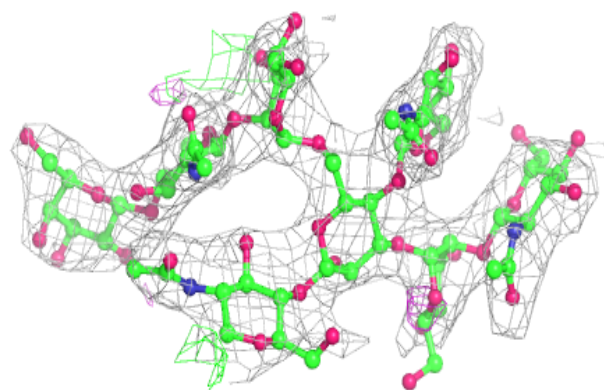


Electron density around Chain L:

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

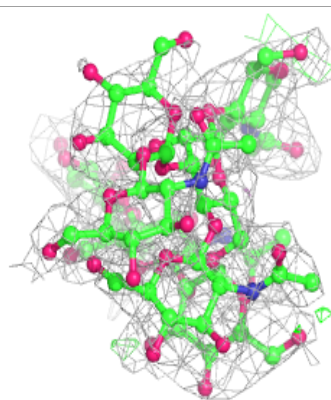
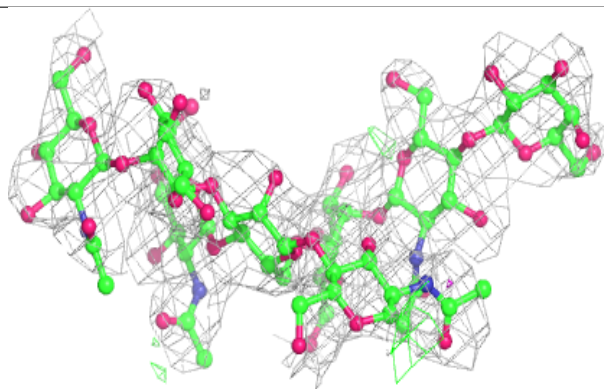
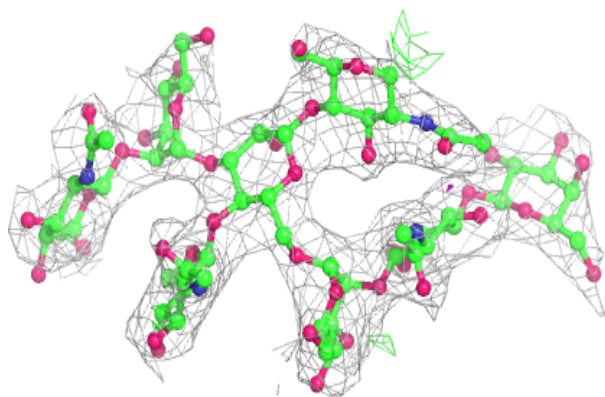
**Electron density around Chain M:**

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

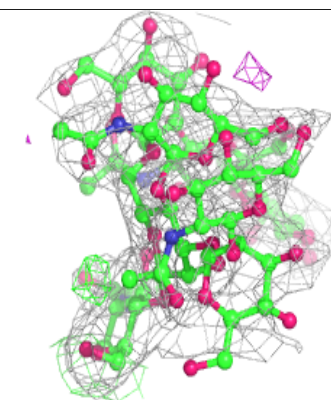
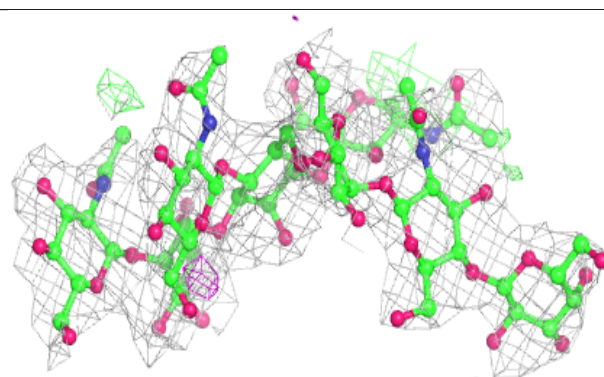
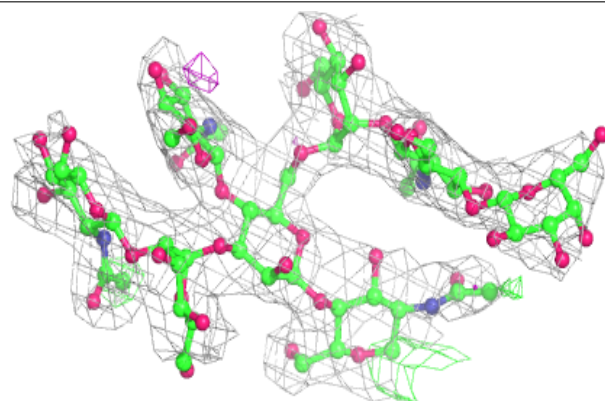


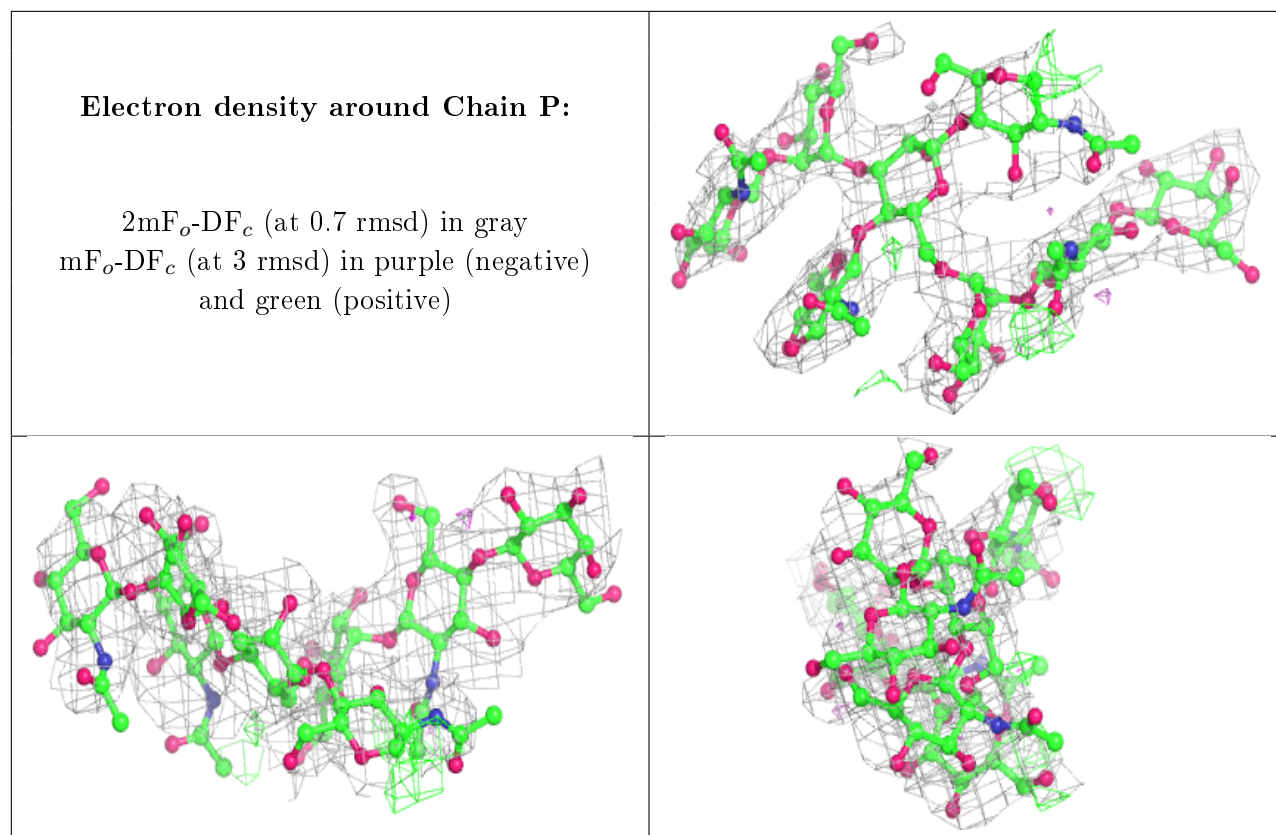
Electron density around Chain N:

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

**Electron density around Chain O:**

$2mF_o-DF_c$ (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
4	CA	H	2002	1/1	0.86	0.08	72,72,72,72	0
4	CA	G	2002	1/1	0.87	0.09	61,61,61,61	0
4	CA	D	2002	1/1	0.90	0.08	44,44,44,44	0
3	MN	C	2001	1/1	0.95	0.06	41,41,41,41	0
3	MN	H	2001	1/1	0.95	0.04	63,63,63,63	0
4	CA	C	2002	1/1	0.95	0.07	42,42,42,42	0
4	CA	E	2002	1/1	0.95	0.12	40,40,40,40	0
3	MN	D	2001	1/1	0.96	0.04	40,40,40,40	0
4	CA	A	2002	1/1	0.96	0.06	34,34,34,34	0
4	CA	B	2002	1/1	0.96	0.07	35,35,35,35	0
3	MN	A	2001	1/1	0.97	0.05	33,33,33,33	0
3	MN	G	2001	1/1	0.97	0.07	54,54,54,54	0
4	CA	F	2002	1/1	0.97	0.05	44,44,44,44	0
3	MN	E	2001	1/1	0.97	0.06	40,40,40,40	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
3	MN	B	2001	1/1	0.98	0.04	34,34,34,34	0
3	MN	F	2001	1/1	0.99	0.04	43,43,43,43	0

6.5 Other polymers [i](#)

There are no such residues in this entry.