



Full wwPDB X-ray Structure Validation Report ⓘ

Aug 6, 2020 – 05:07 PM BST

PDB ID : 1GW0
Title : Crystal Structure of Laccase from *Melanocarpus albomyces* in Four Copper Form
Authors : Hakulinen, N.; Kiiskinen, L.-L.; Kruus, K.; Saloheimo, M.; Koivula, A.; Rouvinen, J.
Deposited on : 2002-03-01
Resolution : 2.40 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

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

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

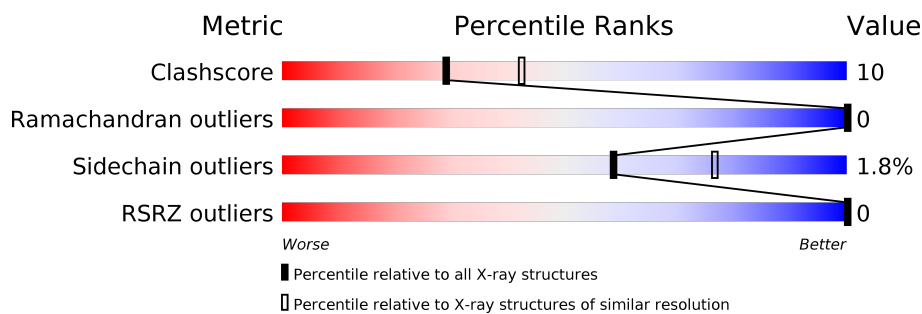
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION




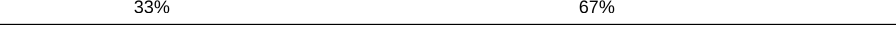
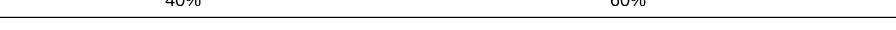
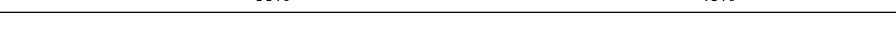
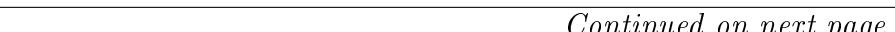
The reported resolution of this entry is 2.40 Å.

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.



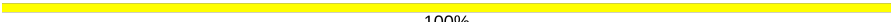


Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
Clashscore	141614	4398 (2.40-2.40)
Ramachandran outliers	138981	4318 (2.40-2.40)
Sidechain outliers	138945	4319 (2.40-2.40)
RSRZ outliers	127900	3811 (2.40-2.40)

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	559	
1	B	559	
2	C	7	
3	D	3	
4	E	5	
4	J	5	
5	F	2	

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Mol	Chain	Length	Quality of chain
5	G	2	 100%
5	I	2	 50% 50%
6	H	6	 67% 33%

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	BMA	C	6	-	-	X	-

2 Entry composition [i](#)

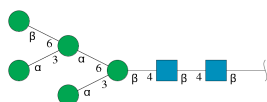
There are 12 unique types of molecules in this entry. The entry contains 10060 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a protein called LACCASE-1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	559	Total	C	N	O	S	0	0	0
			4369	2764	759	831	15			
1	B	559	Total	C	N	O	S	0	0	0
			4369	2764	759	831	15			

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



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	C	7	Total	C	N	O	0	0	0
			83	46	2	35			

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



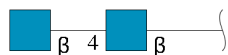
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
3	D	3	Total	C	N	O	0	0	0
			39	22	2	15			

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



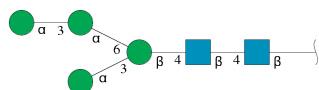
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
4	E	5	Total	C	N	O	0	0	0
			61	34	2	25			
4	J	5	Total	C	N	O	0	0	0
			61	34	2	25			

- Molecule 5 is an oligosaccharide called 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.



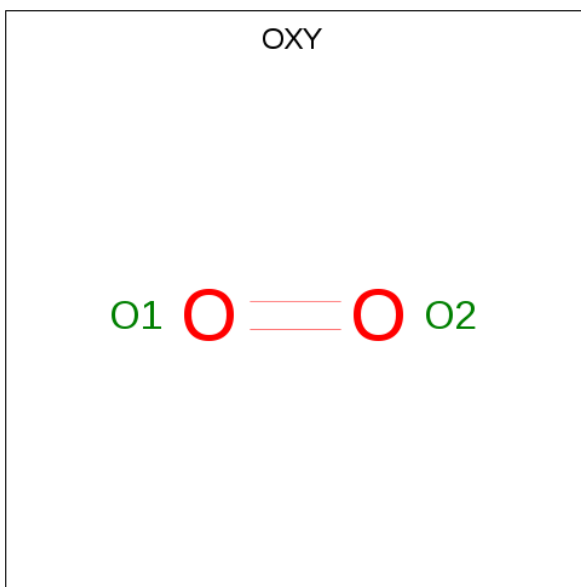
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
5	F	2	Total	C	N	O	0	0	0
			28	16	2	10			
5	G	2	Total	C	N	O	0	0	0
			28	16	2	10			
5	I	2	Total	C	N	O	0	0	0
			28	16	2	10			

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



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
6	H	6	Total	C	N	O	0	0	0
			72	40	2	30			

- Molecule 7 is OXYGEN MOLECULE (three-letter code: OXY) (formula: O₂).



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

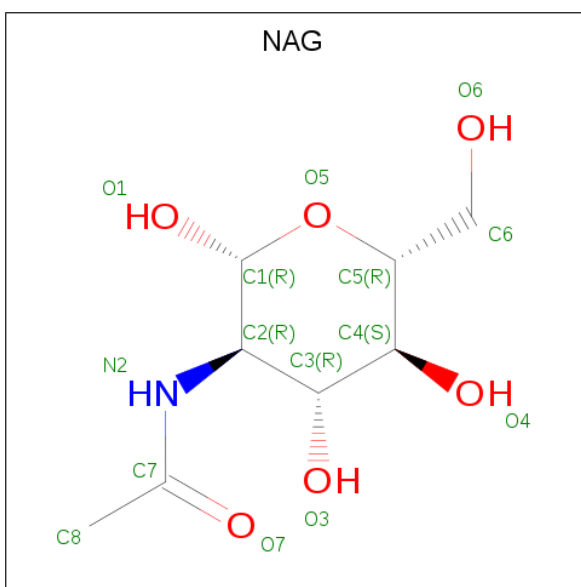
- Molecule 8 is COPPER (II) ION (three-letter code: CU) (formula: Cu).

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

- Molecule 9 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

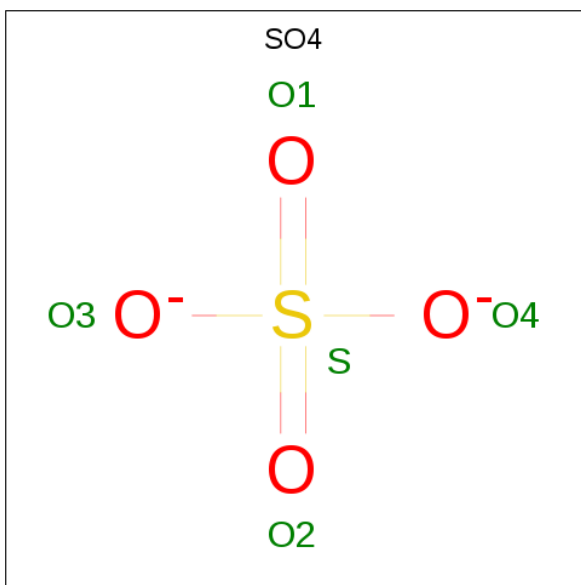
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
9	B	1	Total Cl 1 1	0	0
9	A	1	Total Cl 1 1	0	0

- Molecule 10 is 2-acetamido-2-deoxy-beta-D-glucopyranose (three-letter code: NAG) (formula: C₈H₁₅NO₆).



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

- Molecule 11 is SULFATE ION (three-letter code: SO4) (formula: O₄S).



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

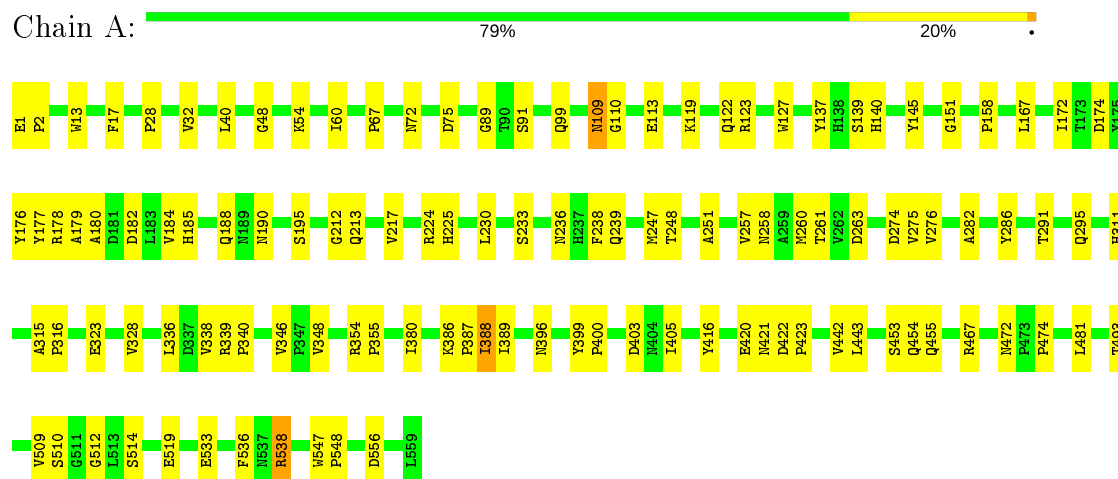
- Molecule 12 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
12	A	408	Total	O	0	0
			408	408		
12	B	387	Total	O	0	0
			387	387		

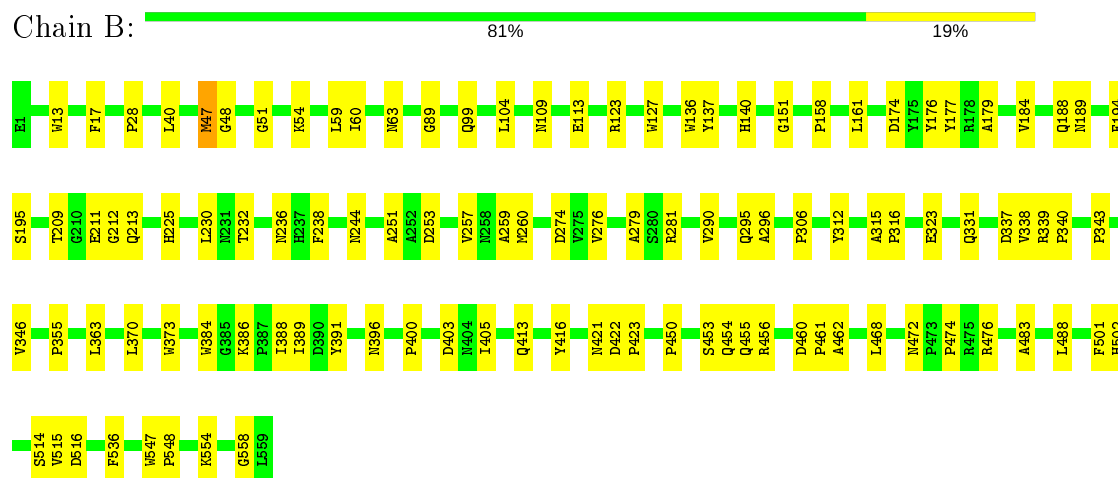
3 Residue-property plots

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

• Molecule 1: LACCASE-1



• Molecule 1: LACCASE-1



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





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

Chain D:



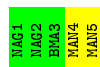
- Molecule 4: alpha-D-mannopyranose-(1-2)-alpha-D-mannopyranose-(1-3)-beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain E:



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

Chain J:



- Molecule 5: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain F:



- Molecule 5: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain G:



- Molecule 5: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain I:



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

Chain H: 

MAN1
MAN2
MAN3
MAN4
MAN5
MAN6

4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	62.51Å 72.30Å 88.94Å 110.40° 95.23° 109.74°	Depositor
Resolution (Å)	100.00 – 2.40 80.99 – 2.40	Depositor EDS
% Data completeness (in resolution range)	95.0 (100.00-2.40) 89.2 (80.99-2.40)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.94 (at 2.40Å)	Xtriage
Refinement program	CNS 1.0	Depositor
R, R_{free}	0.189 , 0.248 0.189 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å ²)	21.1	Xtriage
Anisotropy	0.553	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 48.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	10060	wwPDB-VP
Average B, all atoms (Å ²)	24.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.02% 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: BMA, NAG, CL, OXY, SO4, CU, 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	0.34	0/4506	0.64	1/6191 (0.0%)
1	B	0.34	0/4506	0.63	1/6191 (0.0%)
All	All	0.34	0/9012	0.64	2/12382 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	89	GLY	N-CA-C	-5.81	98.57	113.10
1	B	89	GLY	N-CA-C	-5.02	100.56	113.10

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	4369	0	4111	98	0
1	B	4369	0	4110	75	0
2	C	83	0	70	8	0
3	D	39	0	34	1	0
4	E	61	0	52	1	0
4	J	61	0	52	3	0
5	F	28	0	25	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
5	G	28	0	25	2	0
5	I	28	0	25	0	0
6	H	72	0	61	1	0
7	A	2	0	0	0	0
7	B	2	0	0	0	0
8	A	4	0	0	0	0
8	B	4	0	0	0	0
9	A	1	0	0	0	0
9	B	1	0	0	0	0
10	A	28	0	26	2	0
10	B	70	0	65	2	0
11	A	10	0	0	0	0
11	B	5	0	0	0	0
12	A	408	0	0	8	0
12	B	387	0	0	8	0
All	All	10060	0	8656	172	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 10.

All (172) 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:247:MET:HB3	1:A:275:VAL:HG11	1.39	0.99
1:A:1:GLU:HB2	1:A:2:PRO:HD2	1.47	0.93
1:A:185:HIS:CD2	2:C:6:BMA:H2	2.04	0.92
1:A:400:PRO:HG2	1:A:403:ASP:OD2	1.82	0.79
1:B:453:SER:HB3	1:B:455:GLN:HG3	1.66	0.78
1:A:257:VAL:HG23	1:A:339:ARG:O	1.85	0.76
2:C:4:MAN:H62	2:C:6:BMA:O2	1.85	0.75
1:A:387:PRO:HG2	1:A:536:PHE:HZ	1.53	0.74
1:A:99:GLN:HE22	1:A:127:TRP:HB3	1.52	0.73
1:B:48:GLY:HA3	12:B:717:HOH:O	1.87	0.72
1:A:13:TRP:HB2	1:A:158:PRO:HG3	1.71	0.71
1:B:13:TRP:HB2	1:B:158:PRO:HG3	1.73	0.70
1:A:339:ARG:HH21	1:A:474:PRO:HD3	1.56	0.69
1:B:281:ARG:NH2	10:B:624:NAG:H82	2.08	0.69
1:A:323:GLU:OE1	4:E:2:NAG:H61	1.93	0.69
1:A:251:ALA:HB3	1:A:274:ASP:HB2	1.76	0.68
1:B:338:VAL:HG23	1:B:474:PRO:HB3	1.75	0.68
1:A:190:ASN:HD21	2:C:6:BMA:H5	1.58	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:212:GLY:O	1:B:213:GLN:HG2	1.95	0.67
1:A:354:ARG:HH11	1:A:354:ARG:HB3	1.60	0.66
1:A:354:ARG:NH1	1:A:354:ARG:HB3	2.11	0.65
1:B:386:LYS:HE3	1:B:391:TYR:CZ	2.31	0.65
1:B:453:SER:O	1:B:454:GLN:HB2	1.96	0.65
1:A:190:ASN:ND2	2:C:6:BMA:H5	2.13	0.64
1:A:167:LEU:O	10:A:627:NAG:H83	1.98	0.64
1:B:184:VAL:O	1:B:188:GLN:HG3	1.98	0.63
1:A:54:LYS:HD2	1:A:177:TYR:HA	1.81	0.63
1:B:405:ILE:HD13	1:B:516:ASP:HB3	1.81	0.62
1:A:295:GLN:HA	1:B:295:GLN:OE1	1.99	0.62
1:A:1:GLU:CB	1:A:2:PRO:HD2	2.27	0.61
1:B:54:LYS:HD2	1:B:177:TYR:HA	1.82	0.61
1:B:176:TYR:CE2	1:B:195:SER:HA	2.35	0.61
1:B:59:LEU:HD22	1:B:63:ASN:O	2.01	0.61
1:A:354:ARG:CB	1:A:354:ARG:HH11	2.14	0.61
1:A:247:MET:HB3	1:A:275:VAL:CG1	2.25	0.60
1:B:260:MET:HE1	1:B:337:ASP:C	2.22	0.60
1:A:185:HIS:NE2	2:C:6:BMA:H2	2.17	0.59
1:B:257:VAL:HG12	1:B:340:PRO:HA	1.83	0.59
1:B:47:MET:HE1	1:B:51:GLY:O	2.03	0.58
1:A:399:TYR:HB2	12:A:1005:HOH:O	2.03	0.58
1:B:137:TYR:CZ	1:B:151:GLY:HA3	2.39	0.58
1:B:260:MET:HE1	1:B:338:VAL:N	2.18	0.58
1:A:1:GLU:HB2	1:A:2:PRO:CD	2.28	0.57
1:B:413:GLN:HG3	12:B:1043:HOH:O	2.04	0.57
1:A:48:GLY:HA3	12:A:709:HOH:O	2.04	0.57
1:B:281:ARG:HH22	10:B:624:NAG:H82	1.70	0.56
1:B:461:PRO:HD2	12:B:839:HOH:O	2.05	0.56
1:A:91:SER:O	1:A:139:SER:HA	2.06	0.56
1:B:161:LEU:HD11	1:B:259:ALA:HB2	1.88	0.56
1:B:209:THR:OG1	1:B:211:GLU:HB2	2.06	0.56
1:B:355:PRO:HD2	12:B:895:HOH:O	2.05	0.55
1:A:420:GLU:HG2	12:A:743:HOH:O	2.06	0.55
1:A:67:PRO:HB3	12:A:1054:HOH:O	2.07	0.55
1:A:257:VAL:HG23	1:A:339:ARG:C	2.26	0.55
1:B:174:ASP:HB3	1:B:236:ASN:HB2	1.89	0.55
1:A:547:TRP:HB3	1:A:548:PRO:HD3	1.88	0.54
1:B:400:PRO:HG2	1:B:403:ASP:OD2	2.07	0.54
1:A:311:HIS:CD2	3:D:1:NAG:H5	2.42	0.54
1:A:387:PRO:HG2	1:A:536:PHE:CZ	2.39	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:236:ASN:HB3	1:A:238:PHE:CE2	2.43	0.54
1:A:40:LEU:CD2	1:A:60:ILE:HG12	2.38	0.54
1:A:110:GLY:O	1:A:387:PRO:HA	2.08	0.54
1:A:185:HIS:HD2	2:C:6:BMA:H2	1.69	0.54
1:A:263:ASP:OD1	4:J:5:MAN:H4	2.08	0.54
1:B:290:VAL:HB	1:B:306:PRO:HG2	1.90	0.54
1:A:178:ARG:HH21	1:A:182:ASP:CG	2.11	0.53
1:A:184:VAL:O	1:A:188:GLN:HG3	2.07	0.53
1:A:212:GLY:O	1:A:213:GLN:HG3	2.07	0.53
1:B:450:PRO:HB2	1:B:453:SER:OG	2.08	0.53
1:A:190:ASN:HD21	2:C:6:BMA:C5	2.20	0.53
1:A:348:VAL:HG12	1:A:443:LEU:HD13	1.91	0.53
1:B:113:GLU:HA	1:B:536:PHE:CE1	2.44	0.53
1:A:137:TYR:CZ	1:A:151:GLY:HA3	2.44	0.53
1:A:422:ASP:N	1:A:423:PRO:HD3	2.24	0.52
1:B:47:MET:HE3	1:B:47:MET:HA	1.91	0.52
1:B:421:ASN:ND2	1:B:483:ALA:HA	2.24	0.52
1:B:40:LEU:HD22	1:B:60:ILE:HG12	1.92	0.52
1:A:32:VAL:HB	1:A:75:ASP:OD1	2.10	0.52
1:B:113:GLU:HA	1:B:536:PHE:HE1	1.75	0.52
1:A:2:PRO:HD3	1:A:13:TRP:CZ2	2.45	0.51
1:A:442:VAL:HB	1:A:474:PRO:HG2	1.92	0.51
1:A:510:SER:OG	1:A:556:ASP:OD1	2.26	0.51
1:A:336:LEU:O	1:A:339:ARG:NH2	2.42	0.51
1:A:315:ALA:HB1	1:A:316:PRO:HD2	1.93	0.50
1:B:236:ASN:HB3	1:B:238:PHE:CE2	2.46	0.50
1:A:260:MET:HE1	1:A:338:VAL:N	2.27	0.49
1:A:217:VAL:HG22	10:A:627:NAG:O7	2.12	0.49
1:A:72:ASN:HB2	1:A:75:ASP:OD2	2.13	0.49
1:B:384:TRP:CZ3	1:B:554:LYS:HD2	2.47	0.49
1:A:311:HIS:HE1	1:A:315:ALA:O	1.96	0.48
1:B:40:LEU:CD2	1:B:60:ILE:HG12	2.43	0.48
1:A:388:ILE:HG12	1:A:405:ILE:HD11	1.95	0.48
1:A:174:ASP:HB3	1:A:236:ASN:HB2	1.96	0.48
1:B:456:ARG:HG3	1:B:456:ARG:HH11	1.77	0.48
1:B:253:ASP:OD2	1:B:476:ARG:HB2	2.14	0.48
1:A:340:PRO:HG2	1:A:472:ASN:OD1	2.13	0.48
1:A:396:ASN:ND2	5:G:1:NAG:C7	2.76	0.48
1:A:257:VAL:HG22	1:A:258:ASN:N	2.28	0.48
1:A:346:VAL:HG22	1:A:416:TYR:OH	2.14	0.48
1:A:453:SER:OG	1:A:455:GLN:HG3	2.14	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:493:THR:HA	1:A:519:GLU:OE2	2.14	0.47
1:A:328:VAL:CG2	4:J:5:MAN:H5	2.43	0.47
1:B:161:LEU:HD11	1:B:259:ALA:CB	2.43	0.47
1:B:502:HIS:CB	1:B:514:SER:HB2	2.45	0.47
1:B:502:HIS:HB3	1:B:514:SER:HB2	1.96	0.47
1:A:355:PRO:HD2	12:A:858:HOH:O	2.14	0.47
1:A:179:ALA:CB	2:C:2:NAG:H81	2.44	0.47
1:B:179:ALA:CB	6:H:2:NAG:H81	2.44	0.47
1:B:547:TRP:HB3	1:B:548:PRO:HD3	1.97	0.47
1:B:340:PRO:HG2	1:B:472:ASN:OD1	2.16	0.46
1:A:338:VAL:CG2	1:A:474:PRO:HB3	2.45	0.46
1:A:386:LYS:HE2	5:G:1:NAG:H61	1.96	0.46
1:A:176:TYR:CE2	1:A:195:SER:HA	2.51	0.46
1:A:282:ALA:HB1	12:A:1049:HOH:O	2.14	0.46
1:B:17:PHE:CE1	1:B:28:PRO:HB3	2.51	0.46
1:B:460:ASP:OD1	1:B:462:ALA:HB3	2.16	0.46
1:B:468:LEU:HD22	1:B:488:LEU:HD23	1.97	0.45
1:A:533:GLU:O	1:A:536:PHE:HB3	2.16	0.45
1:B:386:LYS:HE3	1:B:391:TYR:OH	2.16	0.44
1:A:2:PRO:HB3	1:A:13:TRP:CZ3	2.52	0.44
1:A:328:VAL:HG21	4:J:5:MAN:H5	1.98	0.44
1:A:286:TYR:HA	12:A:867:HOH:O	2.18	0.44
1:A:113:GLU:HA	1:A:536:PHE:CE1	2.52	0.44
1:B:388:ILE:HD11	1:B:403:ASP:CG	2.38	0.44
1:A:328:VAL:HG23	12:B:794:HOH:O	2.17	0.44
1:B:296:ALA:N	12:B:704:HOH:O	2.42	0.43
1:B:99:GLN:HE22	1:B:127:TRP:HB3	1.84	0.43
1:A:239:GLN:HB2	1:A:291:THR:O	2.17	0.43
1:B:422:ASP:N	1:B:423:PRO:HD3	2.33	0.43
1:A:275:VAL:HG12	1:A:276:VAL:N	2.34	0.43
1:B:225:HIS:O	1:B:276:VAL:HA	2.19	0.43
1:B:502:HIS:HB3	1:B:514:SER:CB	2.48	0.43
1:B:391:TYR:HB3	1:B:396:ASN:O	2.18	0.43
1:B:343:PRO:HA	1:B:472:ASN:OD1	2.19	0.43
1:A:453:SER:O	1:A:454:GLN:HB2	2.19	0.43
1:A:91:SER:OG	1:A:109:ASN:HB3	2.19	0.42
1:B:251:ALA:HB3	1:B:274:ASP:HB2	2.01	0.42
1:B:47:MET:HB3	12:B:835:HOH:O	2.19	0.42
1:B:331:GLN:OE1	1:B:450:PRO:HA	2.19	0.42
1:B:501:PHE:HB3	1:B:515:VAL:HG22	2.00	0.42
1:A:380:ILE:HG22	1:A:512:GLY:HA2	2.02	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:501:PHE:HB3	1:B:515:VAL:CG2	2.49	0.42
1:A:122:GLN:O	1:A:538:ARG:NH2	2.43	0.42
1:A:17:PHE:CE1	1:A:28:PRO:HB3	2.54	0.42
1:A:257:VAL:HG21	1:A:338:VAL:HB	2.01	0.42
1:B:136:TRP:CE3	1:B:232:THR:HG22	2.55	0.41
1:A:1:GLU:N	12:A:720:HOH:O	2.52	0.41
1:B:189:ASN:ND2	1:B:370:LEU:H	2.18	0.41
1:B:363:LEU:HD13	1:B:373:TRP:CE2	2.54	0.41
1:A:172:ILE:HG22	1:A:238:PHE:CE1	2.55	0.41
1:A:380:ILE:HG21	1:A:514:SER:HB3	2.02	0.41
1:A:467:ARG:HG2	1:A:467:ARG:HH11	1.85	0.41
1:B:279:ALA:HA	1:B:312:TYR:OH	2.21	0.41
1:A:453:SER:HA	1:B:194:PHE:CD1	2.55	0.41
1:B:389:ILE:HD13	1:B:389:ILE:HA	1.81	0.41
1:A:339:ARG:NH2	1:A:474:PRO:HD3	2.31	0.41
1:B:338:VAL:CG2	1:B:474:PRO:HB3	2.44	0.41
1:A:509:VAL:HG13	1:A:510:SER:N	2.36	0.41
1:B:315:ALA:HB1	1:B:316:PRO:HD2	2.03	0.41
1:A:389:ILE:HA	1:A:389:ILE:HD13	1.87	0.41
1:B:257:VAL:HA	1:B:340:PRO:HA	2.02	0.41
1:B:346:VAL:HG22	1:B:416:TYR:OH	2.21	0.41
1:A:251:ALA:HA	1:A:257:VAL:HG12	2.03	0.41
1:B:554:LYS:HE2	1:B:558:GLY:CA	2.52	0.41
1:A:225:HIS:O	1:A:276:VAL:HA	2.20	0.40
1:A:180:ALA:O	1:A:184:VAL:HG23	2.21	0.40
1:A:248:THR:HA	1:A:261:THR:HA	2.03	0.40
1:B:295:GLN:O	1:B:296:ALA:HB3	2.22	0.40
1:A:174:ASP:CG	1:A:233:SER:HB3	2.41	0.40
1:A:13:TRP:HB3	1:A:17:PHE:O	2.22	0.40
1:A:421:ASN:HB2	1:A:481:LEU:HD21	2.03	0.40
1:B:323:GLU:HG3	12:B:984:HOH:O	2.20	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	557/559 (100%)	531 (95%)	26 (5%)	0	100	100
1	B	557/559 (100%)	531 (95%)	26 (5%)	0	100	100
All	All	1114/1118 (100%)	1062 (95%)	52 (5%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	478/478 (100%)	469 (98%)	9 (2%)	57	75
1	B	478/478 (100%)	470 (98%)	8 (2%)	60	78
All	All	956/956 (100%)	939 (98%)	17 (2%)	59	76

All (17) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	109	ASN
1	A	119	LYS
1	A	123	ARG
1	A	140	HIS
1	A	145	TYR
1	A	224	ARG
1	A	230	LEU
1	A	388	ILE
1	A	538	ARG
1	B	47	MET
1	B	104	LEU
1	B	109	ASN
1	B	123	ARG
1	B	140	HIS
1	B	230	LEU

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Mol	Chain	Res	Type
1	B	244	ASN
1	B	339	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (18) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	72	ASN
1	A	99	GLN
1	A	144	GLN
1	A	189	ASN
1	A	190	ASN
1	A	208	ASN
1	A	236	ASN
1	A	244	ASN
1	A	413	GLN
1	A	530	GLN
1	B	20	ASN
1	B	99	GLN
1	B	144	GLN
1	B	189	ASN
1	B	208	ASN
1	B	236	ASN
1	B	455	GLN
1	B	530	GLN

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 ⓘ

32 monosaccharides are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The

Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# $ Z > 2$	Counts	RMSZ	# $ Z > 2$
2	NAG	C	1	1,2	14,14,15	0.51	0	17,19,21	0.68	1 (5%)
2	NAG	C	2	2	14,14,15	0.51	0	17,19,21	0.56	0
2	BMA	C	3	2	11,11,12	0.47	0	15,15,17	0.32	0
2	MAN	C	4	2	11,11,12	0.62	0	15,15,17	0.82	1 (6%)
2	MAN	C	5	2	11,11,12	0.48	0	15,15,17	0.55	0
2	BMA	C	6	2	11,11,12	0.51	0	15,15,17	0.48	0
2	MAN	C	7	2	11,11,12	0.53	0	15,15,17	0.67	1 (6%)
3	NAG	D	1	1,3	14,14,15	0.50	0	17,19,21	0.66	0
3	NAG	D	2	3	14,14,15	0.50	0	17,19,21	0.69	1 (5%)
3	BMA	D	3	3	11,11,12	0.48	0	15,15,17	0.31	0
4	NAG	E	1	1,4	14,14,15	0.52	0	17,19,21	0.67	1 (5%)
4	NAG	E	2	4	14,14,15	0.49	0	17,19,21	0.65	0
4	BMA	E	3	4	11,11,12	0.47	0	15,15,17	0.39	0
4	MAN	E	4	4	11,11,12	0.60	0	15,15,17	0.68	1 (6%)
4	MAN	E	5	4	11,11,12	0.50	0	15,15,17	0.50	0
5	NAG	F	1	1,5	14,14,15	0.54	0	17,19,21	0.91	1 (5%)
5	NAG	F	2	5	14,14,15	0.53	0	17,19,21	0.80	1 (5%)
5	NAG	G	1	1,5	14,14,15	0.57	0	17,19,21	0.60	0
5	NAG	G	2	5	14,14,15	0.50	0	17,19,21	0.88	1 (5%)
6	NAG	H	1	1,6	14,14,15	0.48	0	17,19,21	0.66	0
6	NAG	H	2	6	14,14,15	0.49	0	17,19,21	0.64	0
6	BMA	H	3	6	11,11,12	0.49	0	15,15,17	0.33	0
6	MAN	H	4	6	11,11,12	0.55	0	15,15,17	0.67	1 (6%)
6	MAN	H	5	6	11,11,12	0.49	0	15,15,17	0.54	0
6	MAN	H	6	6	11,11,12	0.58	0	15,15,17	0.51	0
5	NAG	I	1	1,5	14,14,15	0.54	0	17,19,21	0.70	0
5	NAG	I	2	5	14,14,15	0.54	0	17,19,21	0.70	1 (5%)
4	NAG	J	1	1,4	14,14,15	0.55	0	17,19,21	0.56	0
4	NAG	J	2	4	14,14,15	0.51	0	17,19,21	0.65	0
4	BMA	J	3	4	11,11,12	0.45	0	15,15,17	0.25	0
4	MAN	J	4	4	11,11,12	0.61	0	15,15,17	0.70	1 (6%)
4	MAN	J	5	4	11,11,12	0.48	0	15,15,17	0.49	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	C	1	1,2	-	0/6/23/26	0/1/1/1
2	NAG	C	2	2	-	0/6/23/26	0/1/1/1
2	BMA	C	3	2	-	0/2/19/22	0/1/1/1
2	MAN	C	4	2	-	2/2/19/22	0/1/1/1
2	MAN	C	5	2	-	0/2/19/22	0/1/1/1
2	BMA	C	6	2	-	0/2/19/22	0/1/1/1
2	MAN	C	7	2	-	2/2/19/22	0/1/1/1
3	NAG	D	1	1,3	-	0/6/23/26	0/1/1/1
3	NAG	D	2	3	-	4/6/23/26	0/1/1/1
3	BMA	D	3	3	-	0/2/19/22	0/1/1/1
4	NAG	E	1	1,4	-	0/6/23/26	0/1/1/1
4	NAG	E	2	4	-	2/6/23/26	0/1/1/1
4	BMA	E	3	4	-	1/2/19/22	0/1/1/1
4	MAN	E	4	4	-	1/2/19/22	0/1/1/1
4	MAN	E	5	4	-	2/2/19/22	0/1/1/1
5	NAG	F	1	1,5	-	4/6/23/26	0/1/1/1
5	NAG	F	2	5	-	2/6/23/26	0/1/1/1
5	NAG	G	1	1,5	-	5/6/23/26	0/1/1/1
5	NAG	G	2	5	-	4/6/23/26	0/1/1/1
6	NAG	H	1	1,6	-	2/6/23/26	0/1/1/1
6	NAG	H	2	6	-	2/6/23/26	0/1/1/1
6	BMA	H	3	6	-	0/2/19/22	0/1/1/1
6	MAN	H	4	6	-	2/2/19/22	0/1/1/1
6	MAN	H	5	6	-	2/2/19/22	0/1/1/1
6	MAN	H	6	6	-	0/2/19/22	0/1/1/1
5	NAG	I	1	1,5	-	0/6/23/26	0/1/1/1
5	NAG	I	2	5	-	2/6/23/26	0/1/1/1
4	NAG	J	1	1,4	-	0/6/23/26	0/1/1/1
4	NAG	J	2	4	-	1/6/23/26	0/1/1/1
4	BMA	J	3	4	-	2/2/19/22	0/1/1/1
4	MAN	J	4	4	-	1/2/19/22	0/1/1/1
4	MAN	J	5	4	-	2/2/19/22	0/1/1/1

There are no bond length outliers.

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	F	2	NAG	C2-N2-C7	-2.30	119.62	122.90
5	G	2	NAG	C2-N2-C7	-2.30	119.64	122.90
2	C	4	MAN	C1-O5-C5	2.25	115.24	112.19
6	H	4	MAN	C1-O5-C5	2.16	115.12	112.19
2	C	7	MAN	C1-O5-C5	2.14	115.10	112.19
4	E	1	NAG	C2-N2-C7	-2.14	119.85	122.90
5	I	2	NAG	C2-N2-C7	-2.11	119.90	122.90
4	E	4	MAN	C1-O5-C5	2.05	114.97	112.19
2	C	1	NAG	C2-N2-C7	-2.05	119.99	122.90
4	J	4	MAN	C1-O5-C5	2.02	114.93	112.19
3	D	2	NAG	C2-N2-C7	-2.02	120.02	122.90
5	F	1	NAG	C2-N2-C7	-2.01	120.03	122.90

There are no chirality outliers.

All (45) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	G	1	NAG	C1-C2-N2-C7
5	G	1	NAG	C8-C7-N2-C2
5	G	1	NAG	O7-C7-N2-C2
5	F	2	NAG	C8-C7-N2-C2
5	F	2	NAG	O7-C7-N2-C2
4	E	5	MAN	O5-C5-C6-O6
3	D	2	NAG	C8-C7-N2-C2
6	H	5	MAN	O5-C5-C6-O6
4	J	5	MAN	O5-C5-C6-O6
2	C	4	MAN	O5-C5-C6-O6
6	H	4	MAN	O5-C5-C6-O6
4	E	5	MAN	C4-C5-C6-O6
6	H	5	MAN	C4-C5-C6-O6
5	G	1	NAG	C4-C5-C6-O6
6	H	4	MAN	C4-C5-C6-O6
4	J	5	MAN	C4-C5-C6-O6
3	D	2	NAG	O7-C7-N2-C2
2	C	4	MAN	C4-C5-C6-O6
5	G	1	NAG	O5-C5-C6-O6
3	D	2	NAG	O5-C5-C6-O6
5	G	2	NAG	C4-C5-C6-O6
2	C	7	MAN	O5-C5-C6-O6
5	F	1	NAG	C4-C5-C6-O6
2	C	7	MAN	C4-C5-C6-O6
4	E	2	NAG	C4-C5-C6-O6
5	I	2	NAG	O5-C5-C6-O6

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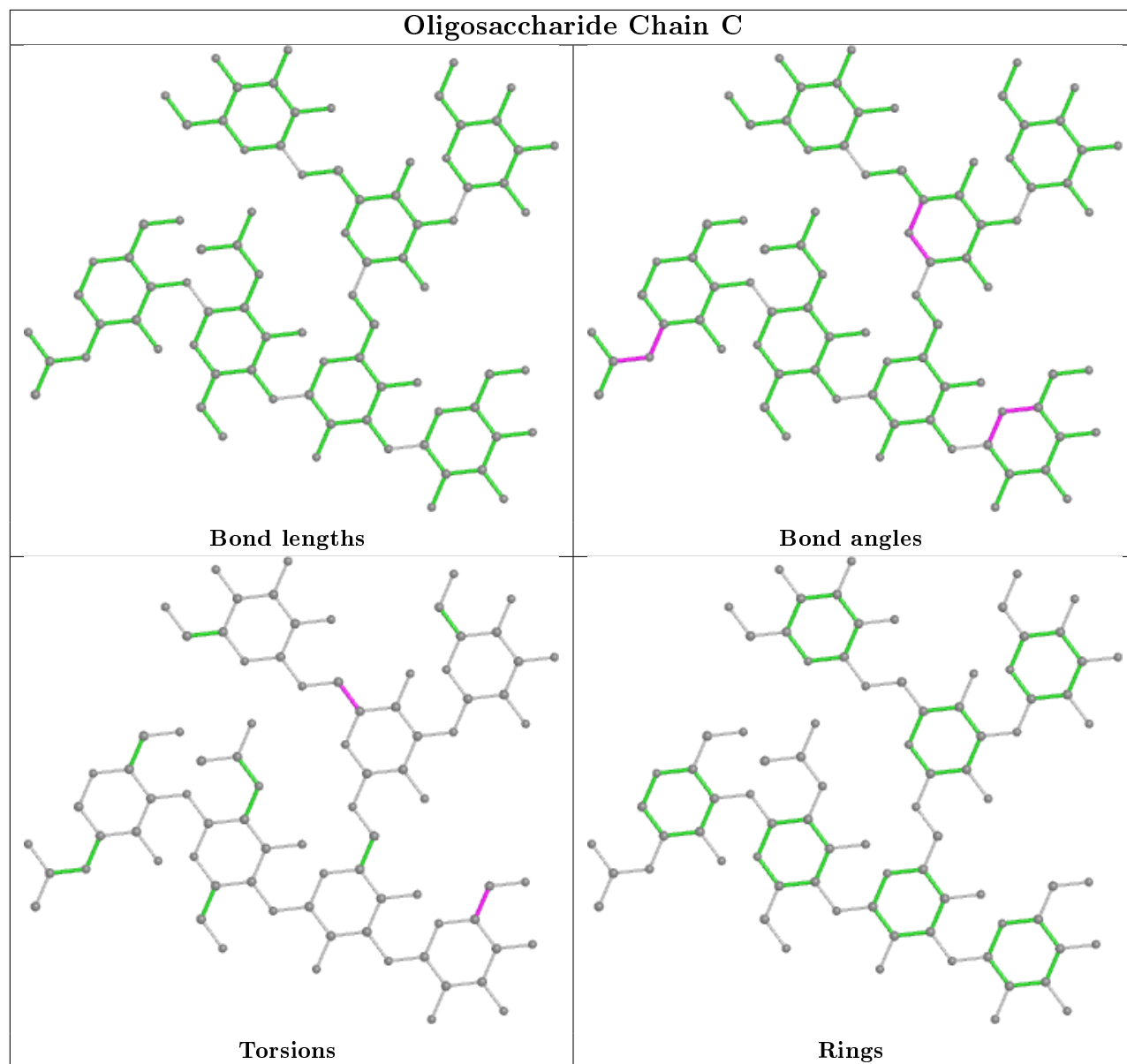
Mol	Chain	Res	Type	Atoms
6	H	1	NAG	C4-C5-C6-O6
5	G	2	NAG	O5-C5-C6-O6
5	F	1	NAG	O5-C5-C6-O6
3	D	2	NAG	C4-C5-C6-O6
4	E	3	BMA	O5-C5-C6-O6
5	G	2	NAG	C8-C7-N2-C2
4	E	2	NAG	O5-C5-C6-O6
6	H	1	NAG	O5-C5-C6-O6
5	F	1	NAG	C8-C7-N2-C2
4	J	3	BMA	C4-C5-C6-O6
4	E	4	MAN	C4-C5-C6-O6
6	H	2	NAG	C4-C5-C6-O6
5	F	1	NAG	O7-C7-N2-C2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

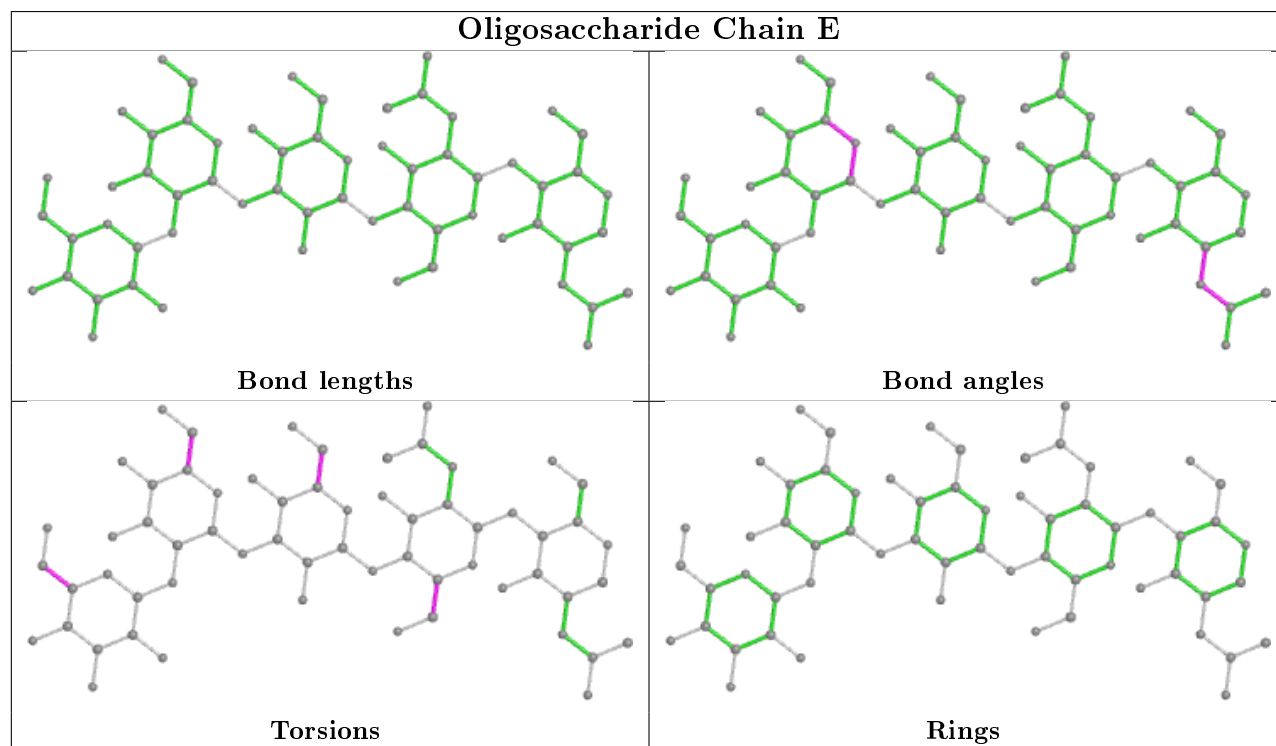
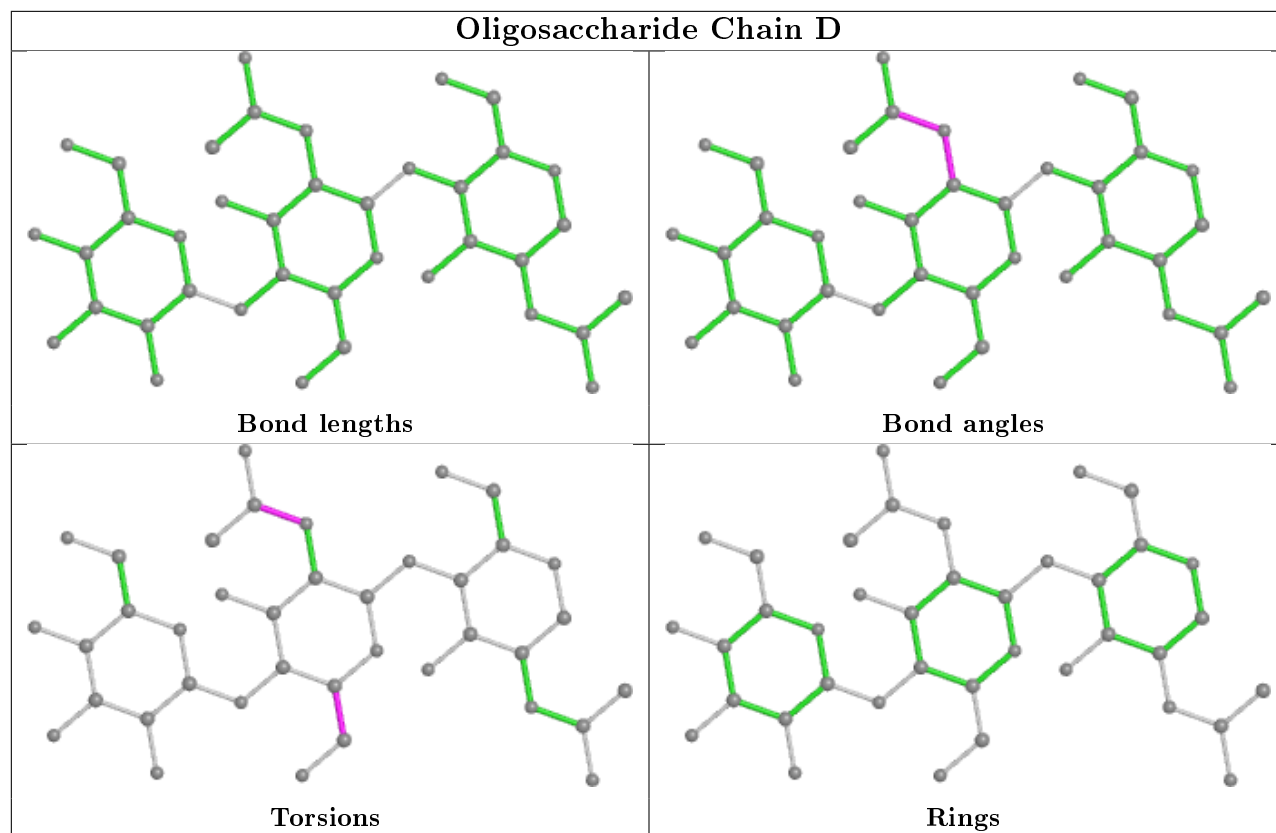
There are no ring outliers.

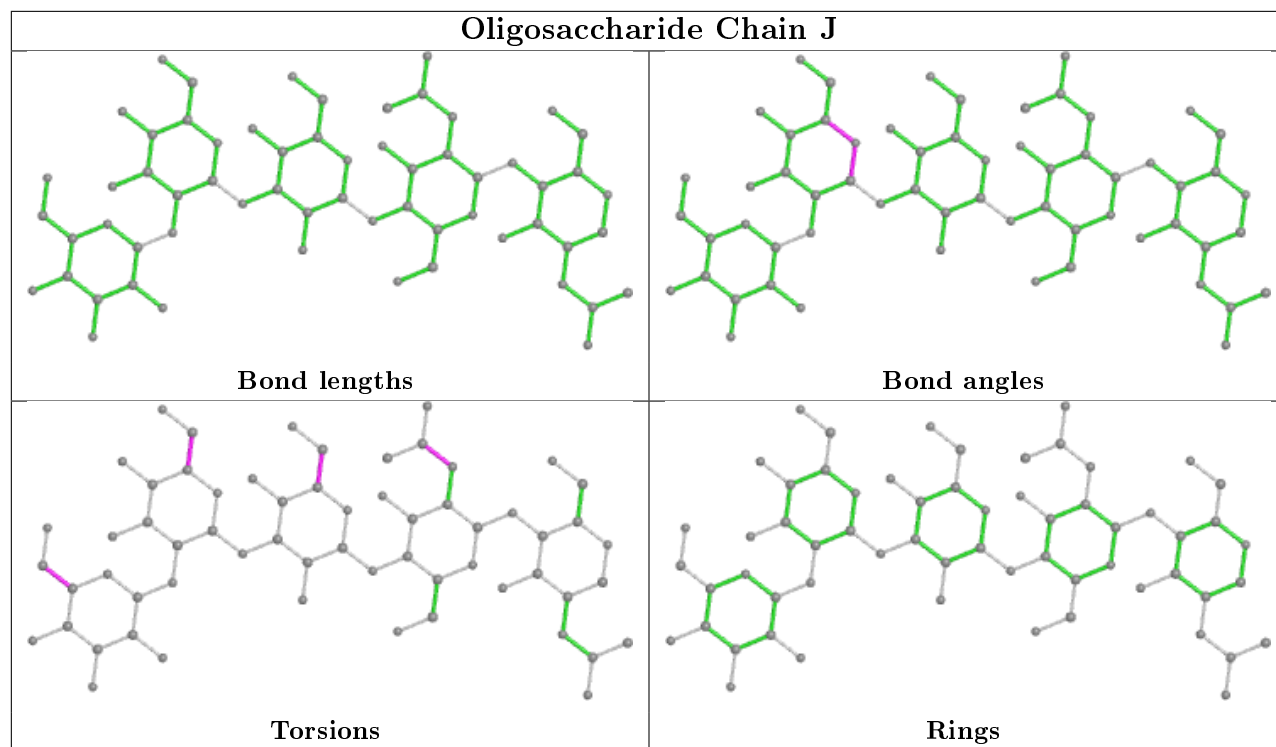
8 monomers are involved in 16 short contacts:

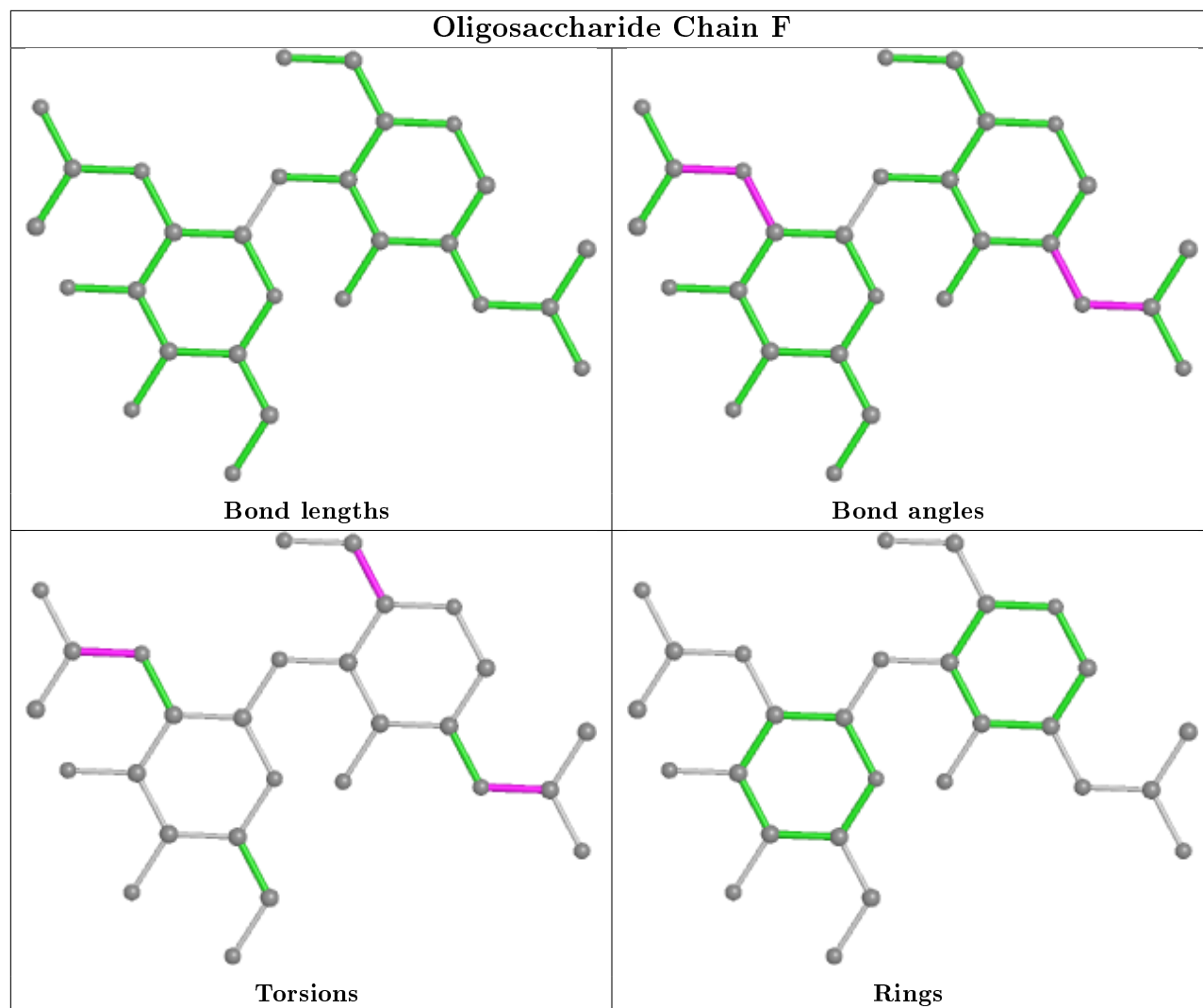
Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	J	5	MAN	3	0
2	C	6	BMA	7	0
2	C	4	MAN	1	0
3	D	1	NAG	1	0
5	G	1	NAG	2	0
2	C	2	NAG	1	0
6	H	2	NAG	1	0
4	E	2	NAG	1	0

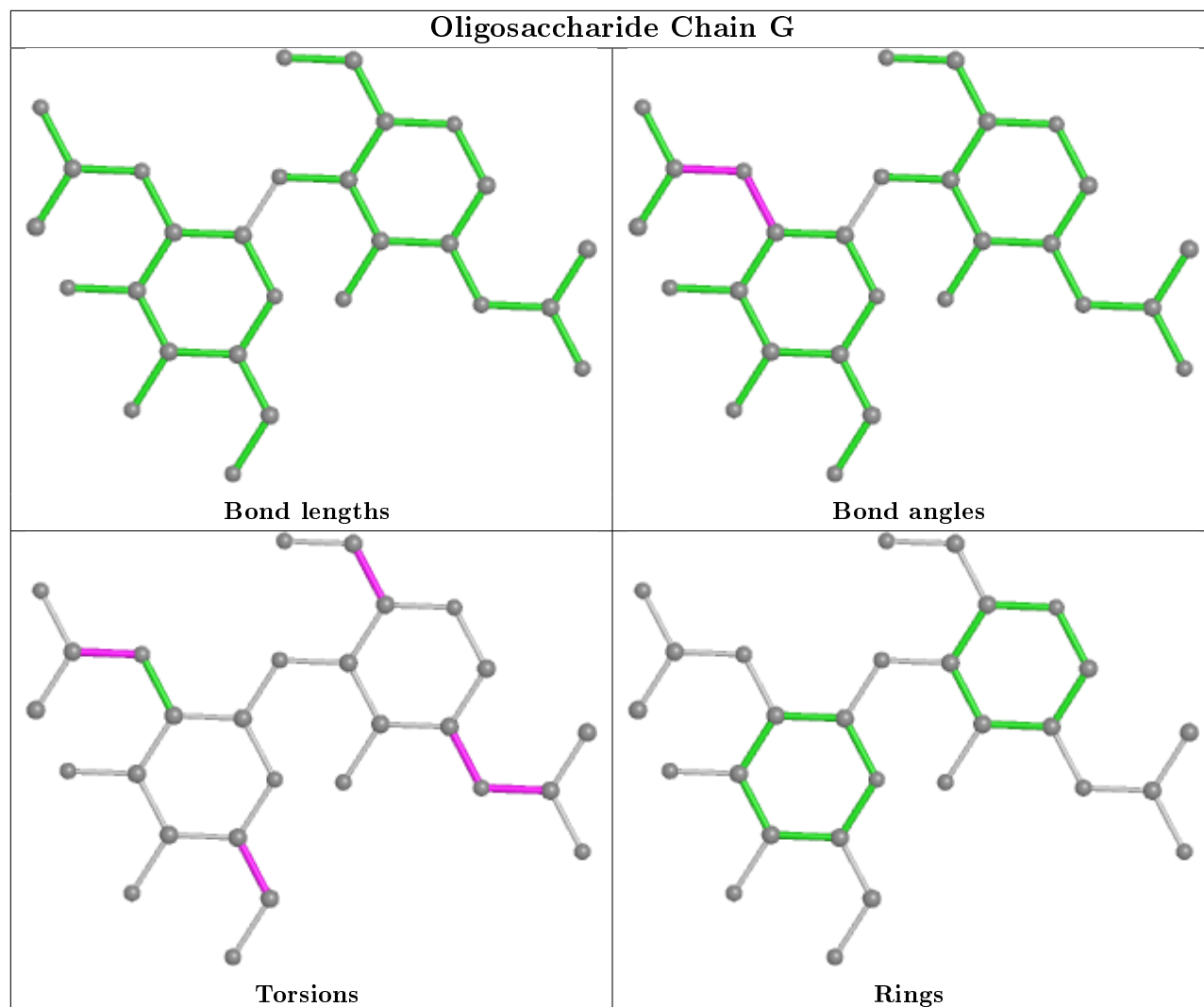
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for oligosaccharide.

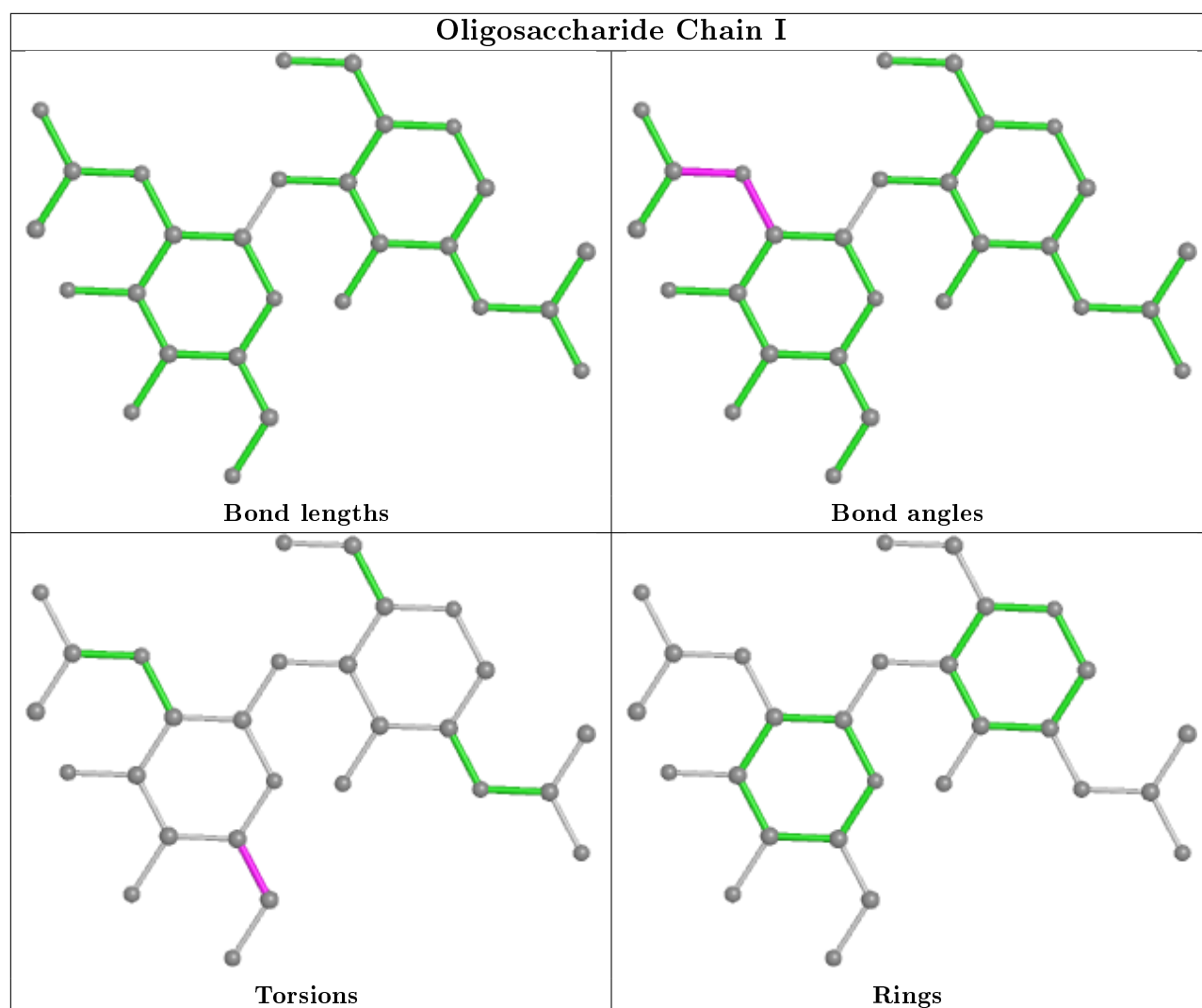


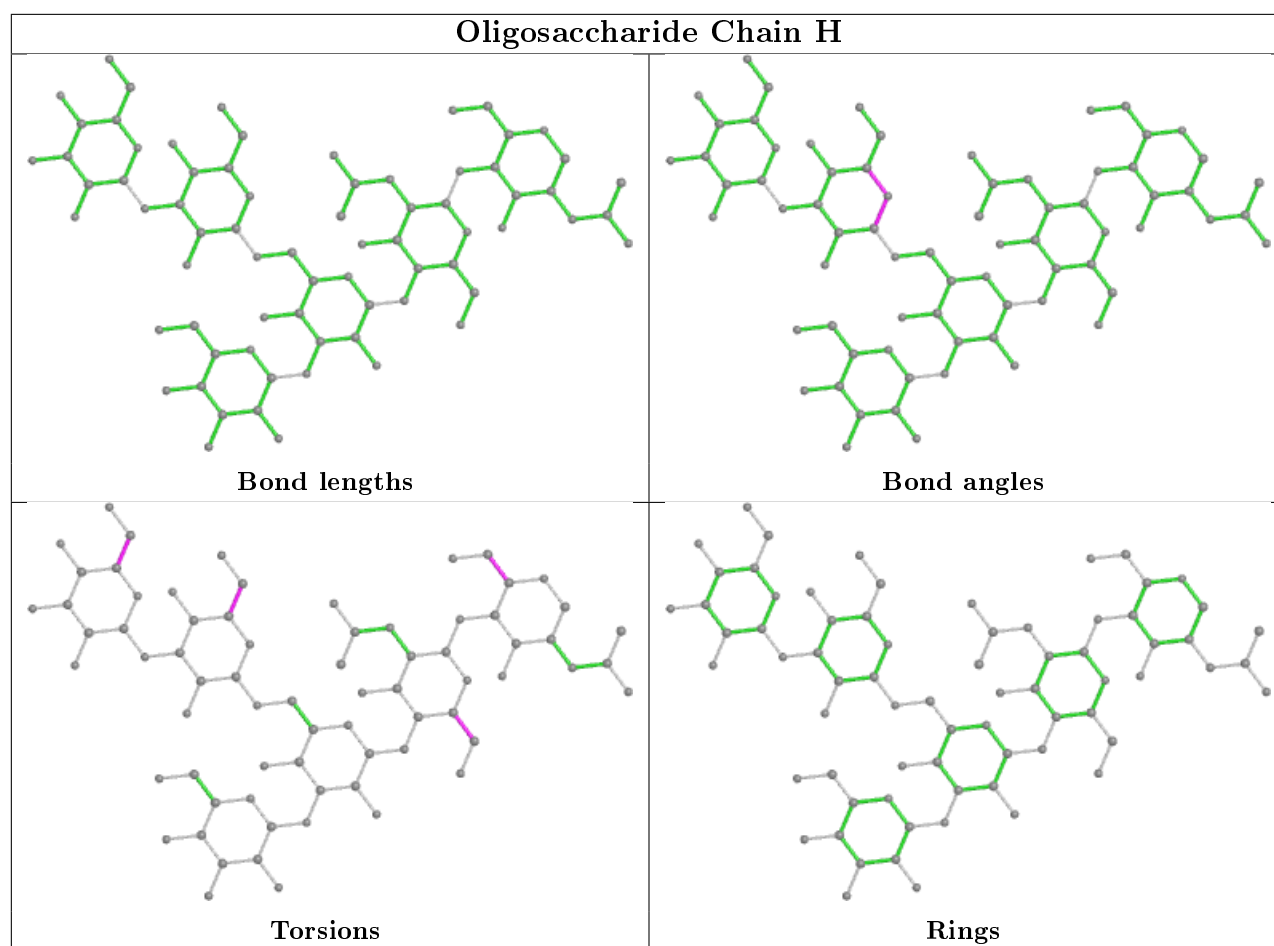












5.6 Ligand geometry [i](#)

Of 22 ligands modelled in this entry, 10 are monoatomic - leaving 12 for Mogul analysis.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
11	SO4	B	625	-	4,4,4	0.27	0	6,6,6	0.07	0
7	OXY	B	601	8	1,1,1	0.10	0	-		
10	NAG	A	627	1	14,14,15	0.51	0	17,19,21	0.83	1 (5%)
11	SO4	A	628	-	4,4,4	0.26	0	6,6,6	0.11	0
10	NAG	A	607	1	14,14,15	0.53	0	17,19,21	0.69	1 (5%)
10	NAG	B	607	1	14,14,15	0.53	0	17,19,21	0.70	1 (5%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
10	NAG	B	621	1	14,14,15	0.50	0	17,19,21	0.63	0
10	NAG	B	622	1	14,14,15	0.56	0	17,19,21	0.64	0
7	OXY	A	601	8	1,1,1	0.09	0	-		
10	NAG	B	623	1	14,14,15	0.65	0	17,19,21	0.76	1 (5%)
11	SO4	A	629	-	4,4,4	0.26	0	6,6,6	0.08	0
10	NAG	B	624	1	14,14,15	0.53	0	17,19,21	0.68	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
10	NAG	A	607	1	-	4/6/23/26	0/1/1/1
10	NAG	B	607	1	-	4/6/23/26	0/1/1/1
10	NAG	B	621	1	-	0/6/23/26	0/1/1/1
10	NAG	B	622	1	-	3/6/23/26	0/1/1/1
10	NAG	A	627	1	-	4/6/23/26	0/1/1/1
10	NAG	B	623	1	-	2/6/23/26	0/1/1/1
10	NAG	B	624	1	-	4/6/23/26	0/1/1/1

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
10	A	627	NAG	C2-N2-C7	-2.43	119.44	122.90
10	A	607	NAG	C2-N2-C7	-2.22	119.74	122.90
10	B	623	NAG	C2-N2-C7	-2.20	119.77	122.90
10	B	607	NAG	C2-N2-C7	-2.17	119.82	122.90

There are no chirality outliers.

All (21) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
10	B	622	NAG	C1-C2-N2-C7
10	B	622	NAG	C8-C7-N2-C2
10	B	622	NAG	O7-C7-N2-C2
10	A	607	NAG	O5-C5-C6-O6
10	A	627	NAG	C8-C7-N2-C2

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Mol	Chain	Res	Type	Atoms
10	A	627	NAG	O7-C7-N2-C2
10	A	607	NAG	C4-C5-C6-O6
10	A	627	NAG	O5-C5-C6-O6
10	B	624	NAG	O5-C5-C6-O6
10	B	607	NAG	C8-C7-N2-C2
10	B	623	NAG	O5-C5-C6-O6
10	A	627	NAG	C4-C5-C6-O6
10	A	607	NAG	C8-C7-N2-C2
10	B	607	NAG	O7-C7-N2-C2
10	B	623	NAG	C4-C5-C6-O6
10	B	607	NAG	O5-C5-C6-O6
10	B	624	NAG	C4-C5-C6-O6
10	B	607	NAG	C4-C5-C6-O6
10	A	607	NAG	O7-C7-N2-C2
10	B	624	NAG	C8-C7-N2-C2
10	B	624	NAG	O7-C7-N2-C2

There are no ring outliers.

2 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
10	A	627	NAG	2	0
10	B	624	NAG	2	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data [i](#)

6.1 Protein, DNA and RNA chains [i](#)

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	559/559 (100%)	-0.77	0 100 100	10, 21, 38, 57	0
1	B	559/559 (100%)	-0.76	0 100 100	7, 22, 37, 58	0
All	All	1118/1118 (100%)	-0.76	0 100 100	7, 21, 37, 58	0

There are no RSRZ outliers to report.

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	E	4	11/12	0.54	0.27	71,72,74,76	0
6	MAN	H	5	11/12	0.59	0.35	94,95,96,96	0
6	MAN	H	4	11/12	0.67	0.34	86,89,92,92	0
4	MAN	J	4	11/12	0.69	0.19	65,67,68,70	0
2	MAN	C	4	11/12	0.70	0.28	72,74,79,82	0
6	MAN	H	6	11/12	0.70	0.31	76,79,80,80	0
5	NAG	F	2	14/15	0.72	0.34	57,60,64,65	0
2	BMA	C	6	11/12	0.72	0.34	84,85,88,89	0
4	MAN	J	5	11/12	0.73	0.23	71,72,74,74	0
4	MAN	E	5	11/12	0.77	0.28	78,79,80,82	0
3	BMA	D	3	11/12	0.80	0.18	63,65,65,66	0

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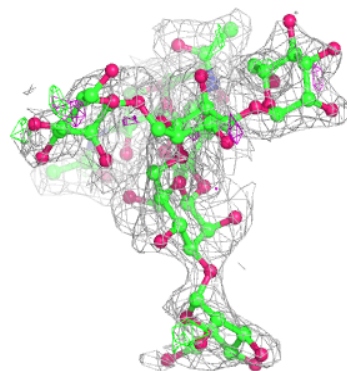
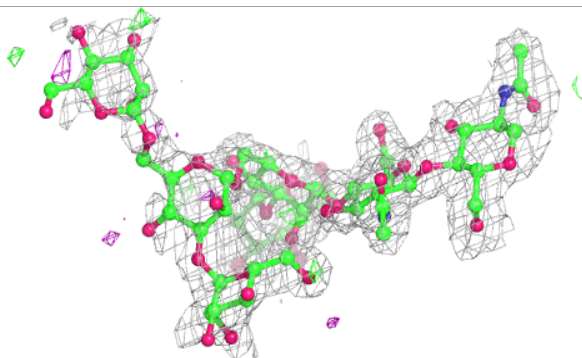
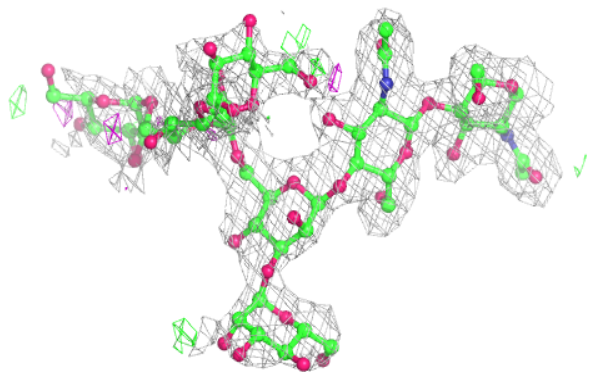
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
2	MAN	C	5	11/12	0.80	0.17	66,70,71,72	0
4	BMA	E	3	11/12	0.82	0.18	57,62,66,68	0
2	MAN	C	7	11/12	0.82	0.28	70,73,74,74	0
6	BMA	H	3	11/12	0.84	0.17	59,67,74,81	0
5	NAG	G	2	14/15	0.84	0.16	48,49,52,54	0
3	NAG	D	2	14/15	0.86	0.18	50,53,57,60	0
2	BMA	C	3	11/12	0.86	0.17	53,59,66,68	0
4	BMA	J	3	11/12	0.87	0.20	53,59,61,63	0
3	NAG	D	1	14/15	0.89	0.14	41,44,47,50	0
5	NAG	I	2	14/15	0.90	0.15	38,40,41,41	0
5	NAG	G	1	14/15	0.90	0.12	39,43,45,48	0
5	NAG	I	1	14/15	0.91	0.14	35,37,37,39	0
5	NAG	F	1	14/15	0.91	0.16	28,36,43,52	0
4	NAG	J	2	14/15	0.91	0.13	36,38,44,51	0
2	NAG	C	2	14/15	0.93	0.12	26,35,38,46	0
4	NAG	J	1	14/15	0.93	0.10	22,27,32,34	0
4	NAG	E	2	14/15	0.94	0.11	36,40,44,50	0
2	NAG	C	1	14/15	0.95	0.10	14,23,26,29	0
4	NAG	E	1	14/15	0.95	0.08	21,25,30,35	0
6	NAG	H	2	14/15	0.95	0.10	24,35,41,50	0
6	NAG	H	1	14/15	0.96	0.09	17,26,29,31	0

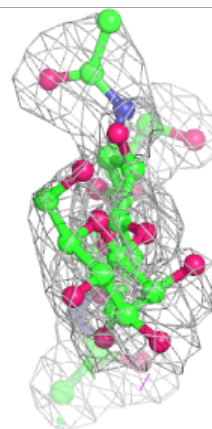
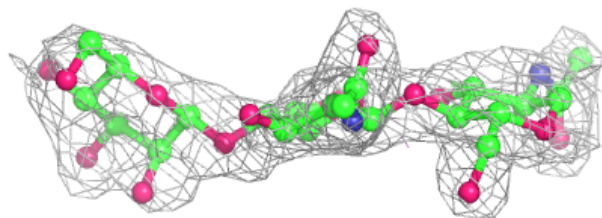
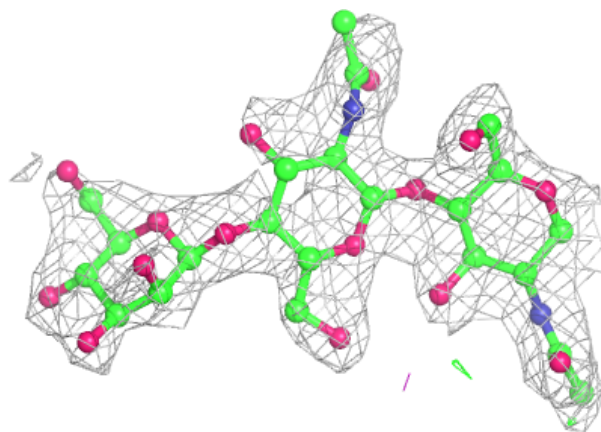
The following is a graphical depiction of the model fit to experimental electron density for oligosaccharide. Each fit is shown from different orientation to approximate a three-dimensional view.

Electron density around Chain C:

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

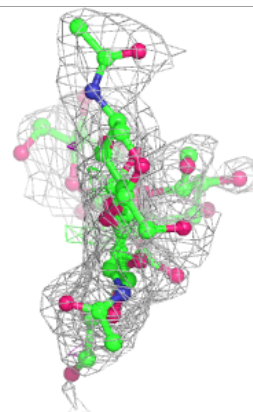
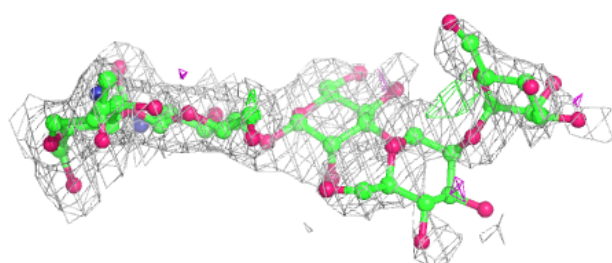
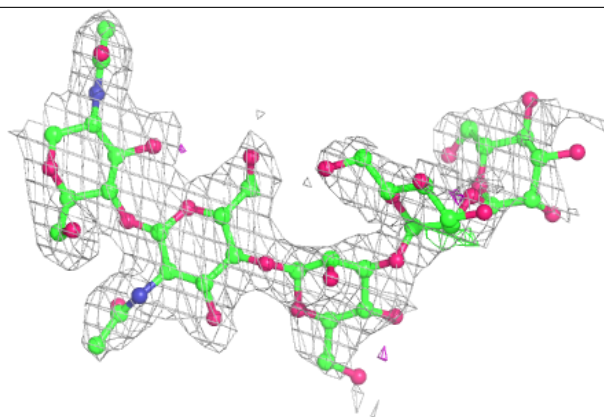
**Electron density around Chain D:**

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

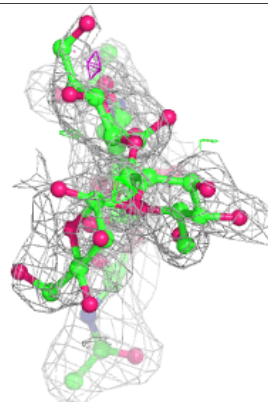
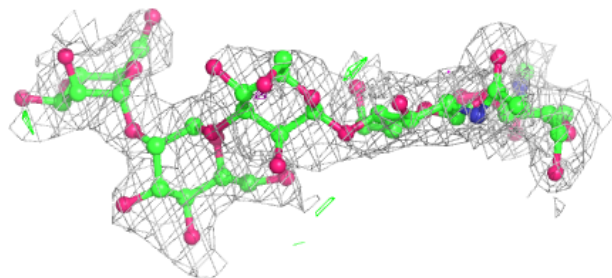
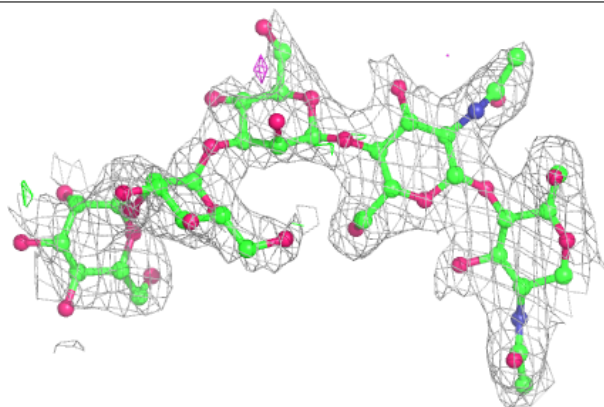


Electron density around Chain E:

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

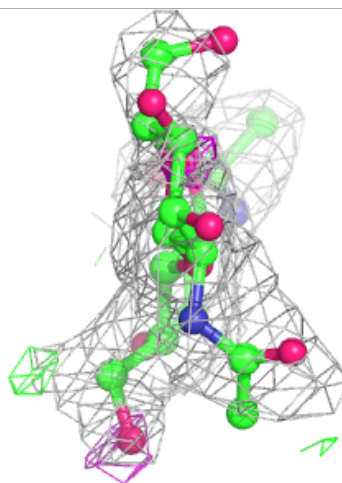
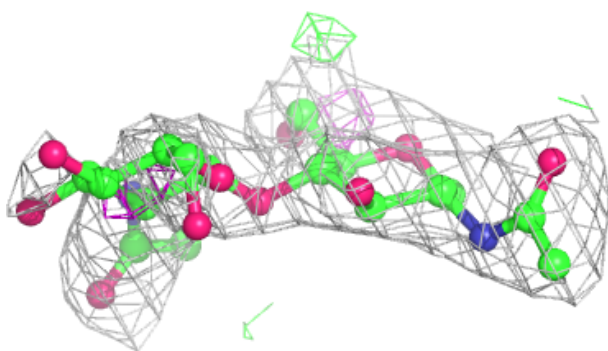
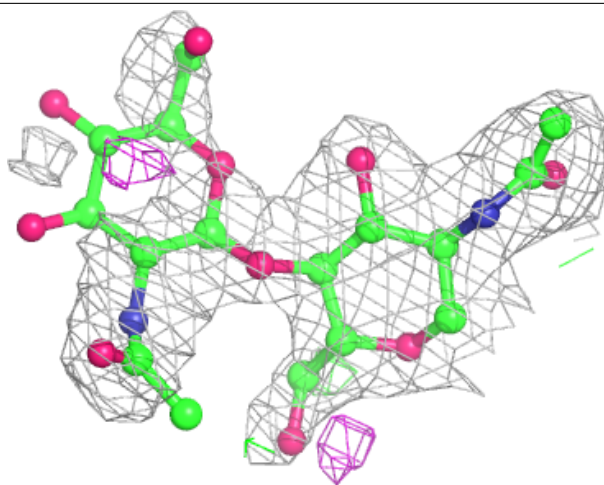
**Electron density around Chain J:**

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



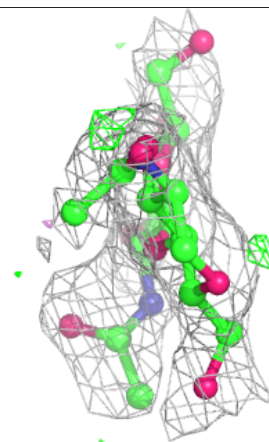
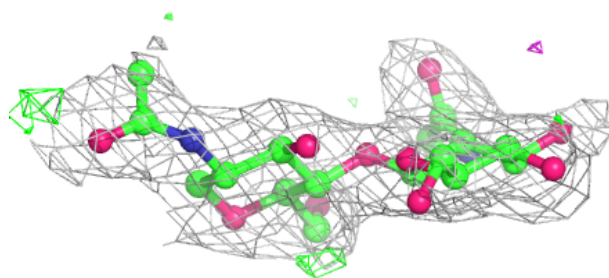
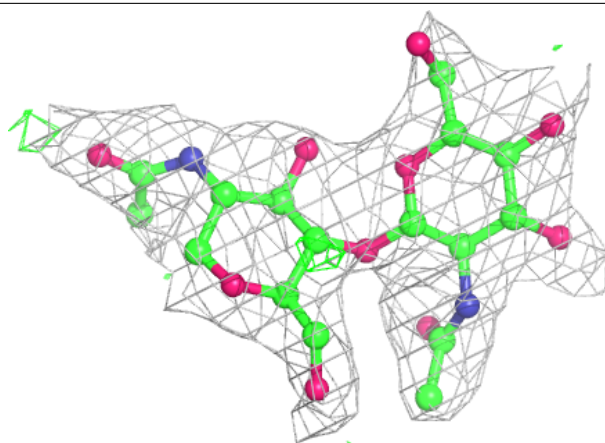
Electron density around Chain F:

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



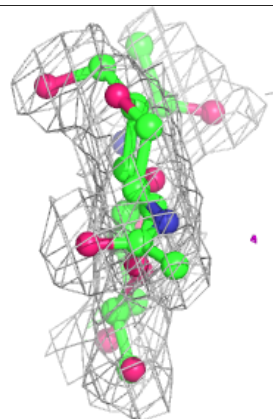
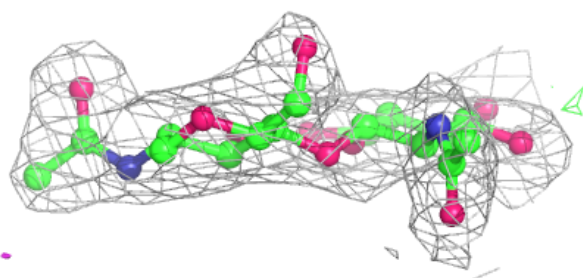
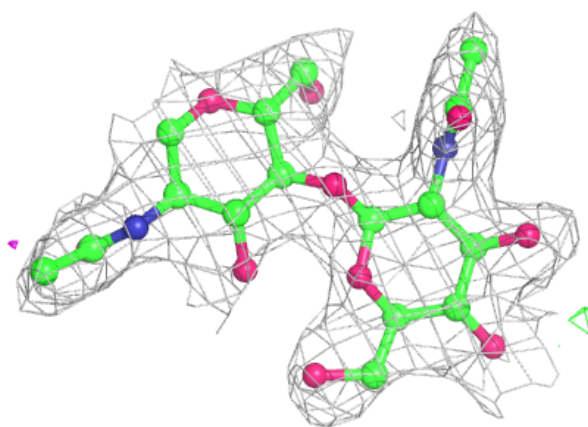
Electron density around Chain G:

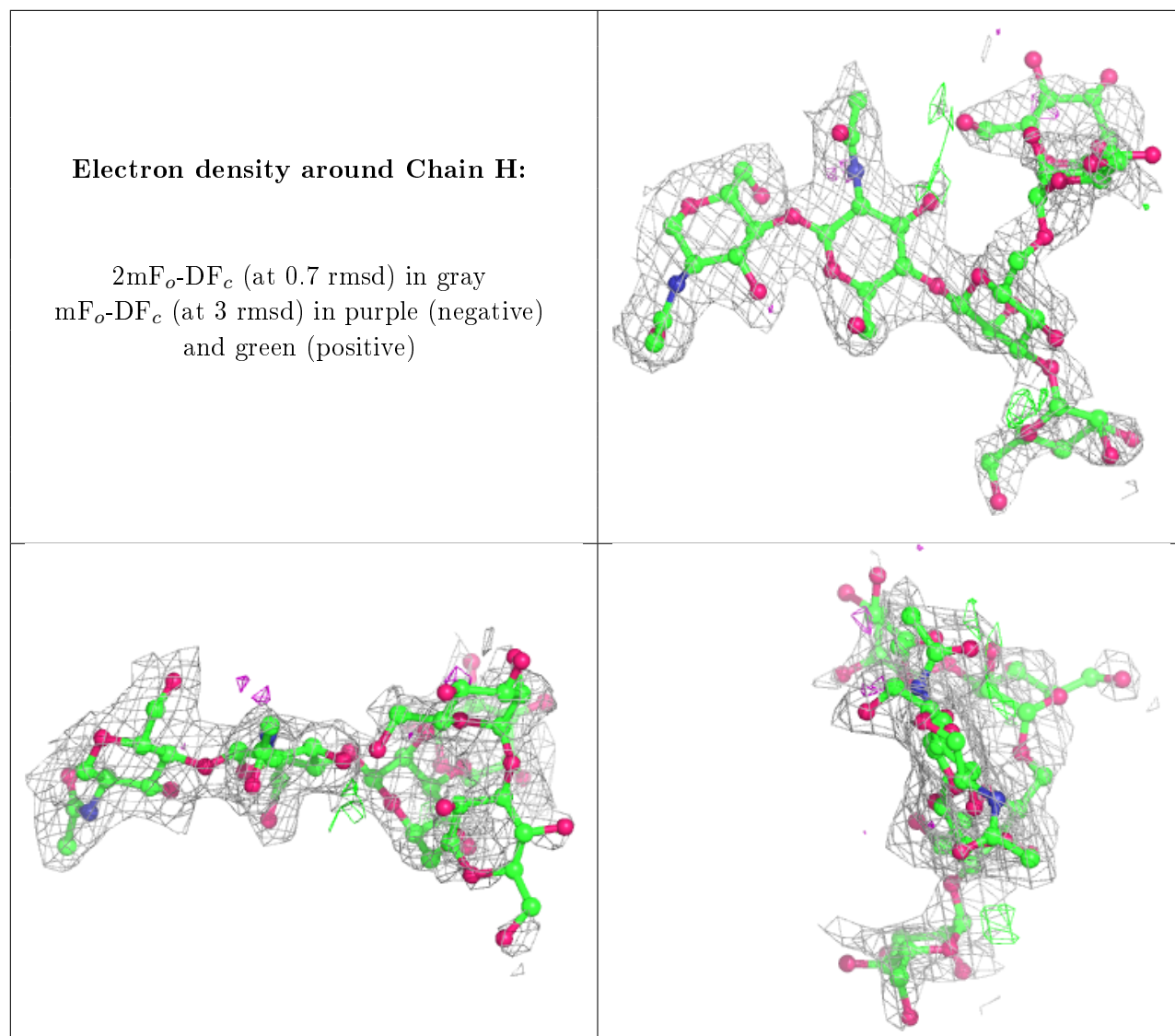
$2mF_o - DF_c$ (at 0.7 rmsd) in gray
 $mF_o - DF_c$ (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around Chain I:

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





6.4 Ligands [i](#)

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
10	NAG	B	624	14/15	0.73	0.21	55,57,60,60	0
10	NAG	B	623	14/15	0.77	0.16	38,44,48,49	0
10	NAG	A	627	14/15	0.82	0.16	46,49,53,55	0
10	NAG	B	622	14/15	0.83	0.18	49,52,54,56	0
11	SO4	A	629	5/5	0.89	0.17	77,77,78,79	0
10	NAG	B	607	14/15	0.89	0.17	34,36,39,40	0
10	NAG	A	607	14/15	0.91	0.13	31,34,37,39	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
10	NAG	B	621	14/15	0.92	0.13	26,30,34,35	0
7	OXY	A	601	2/2	0.94	0.19	22,22,22,25	0
11	SO4	A	628	5/5	0.94	0.10	61,62,62,63	0
11	SO4	B	625	5/5	0.94	0.13	73,74,74,74	0
7	OXY	B	601	2/2	0.97	0.14	19,19,19,23	0
8	CU	A	604	1/1	0.98	0.05	25,25,25,25	0
9	CL	A	606	1/1	0.98	0.07	31,31,31,31	0
8	CU	A	603	1/1	0.99	0.06	23,23,23,23	0
9	CL	B	606	1/1	0.99	0.08	25,25,25,25	0
8	CU	B	604	1/1	0.99	0.06	23,23,23,23	0
8	CU	B	603	1/1	0.99	0.06	28,28,28,28	0
8	CU	A	605	1/1	0.99	0.05	26,26,26,26	0
8	CU	B	605	1/1	0.99	0.07	27,27,27,27	0
8	CU	B	602	1/1	1.00	0.07	25,25,25,25	0
8	CU	A	602	1/1	1.00	0.05	26,26,26,26	0

6.5 Other polymers [i](#)

There are no such residues in this entry.