



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 05:50 PM UTC

PDB ID : 5I8R / pdb_00005i8r
Title : aSMase with zinc
Authors : Zhou, Y.F.; Wei, R.R.
Deposited on : 2016-02-19
Resolution : 3.65 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

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

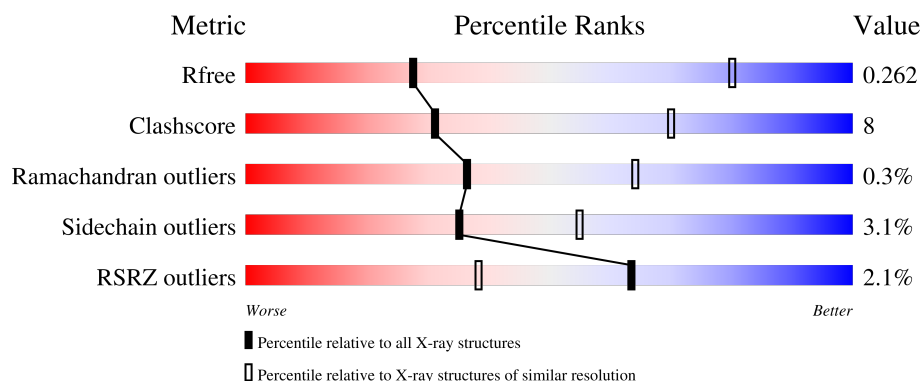
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.65 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



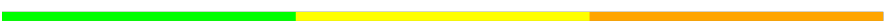
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	1132 (3.76-3.52)
Clashscore	190562	1014 (3.74-3.54)
Ramachandran outliers	187476	1123 (3.76-3.52)
Sidechain outliers	187428	1121 (3.76-3.52)
RSRZ outliers	180081	1130 (3.76-3.52)

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

Mol	Chain	Length	Quality of chain
1	A	583	0% 70% 20% 9%
1	B	583	2% 75% 14% 9%
1	C	583	3% 75% 15% 9%
2	D	2	100%
2	E	2	50% 50%

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Mol	Chain	Length	Quality of chain
2	F	2	 100%
2	G	2	 100%
2	I	2	 100%
3	H	3	 33% 33% 33%

2 Entry composition [i](#)

There are 7 unique types of molecules in this entry. The entry contains 12883 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 Sphingomyelin phosphodiesterase.

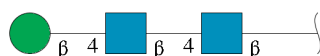
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	528	Total	C	N	O	S	0	0	0
			4161	2684	717	737	23			
1	B	528	Total	C	N	O	S	0	0	0
			4161	2684	717	737	23			
1	C	528	Total	C	N	O	S	0	0	0
			4161	2684	717	737	23			

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



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

- 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	H	3	Total	C	N	O	0	0	0
			39	22	2	15			

- Molecule 4 is 2-acetamido-2-deoxy-beta-D-glucopyranose (CCD ID: NAG) (formula: $C_8H_{15}NO_6$).



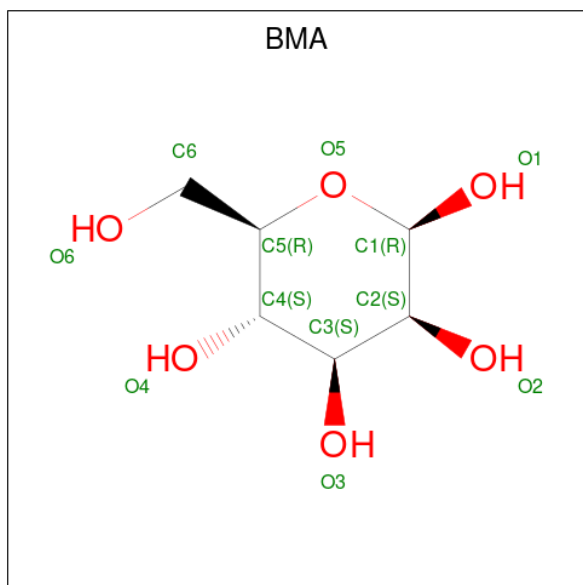
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	1	Total	C	N	O	0	0
			14	8	1	5		
4	A	1	Total	C	N	O	0	0
			14	8	1	5		
4	A	1	Total	C	N	O	0	0
			14	8	1	5		
4	A	1	Total	C	N	O	0	0
			14	8	1	5		
4	A	1	Total	C	N	O	0	0
			14	8	1	5		
4	B	1	Total	C	N	O	0	0
			14	8	1	5		
4	B	1	Total	C	N	O	0	0
			14	8	1	5		
4	B	1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

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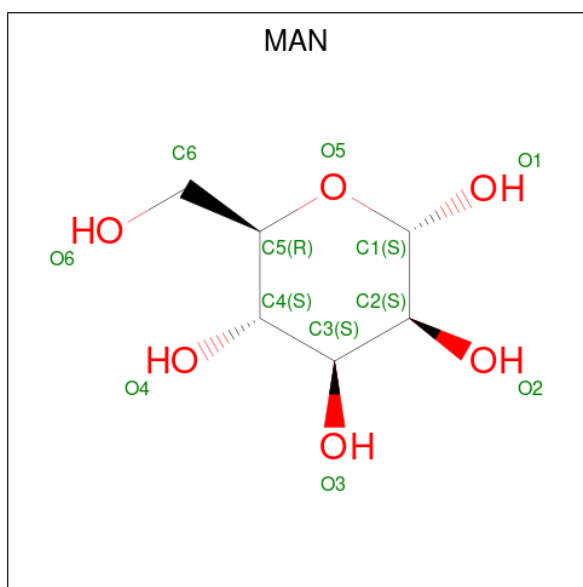
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	C	1	Total	C	N	O	0	0
			14	8	1	5		
4	C	1	Total	C	N	O	0	0
			14	8	1	5		
4	C	1	Total	C	N	O	0	0
			14	8	1	5		

- Molecule 5 is beta-D-mannopyranose (CCD ID: BMA) (formula: $C_6H_{12}O_6$).



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

- Molecule 6 is alpha-D-mannopyranose (CCD ID: MAN) (formula: $C_6H_{12}O_6$).



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

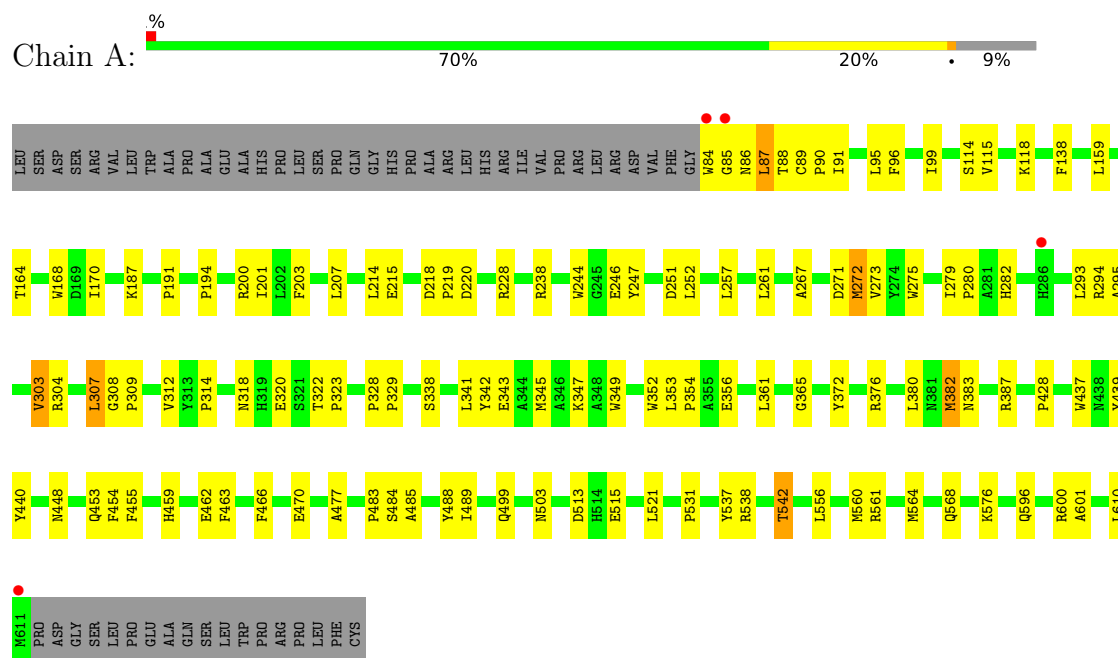
- Molecule 7 is ZINC ION (CCD ID: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	2	Total	Zn	0	0
			2	2		
7	B	2	Total	Zn	0	0
			2	2		
7	C	2	Total	Zn	0	0
			2	2		

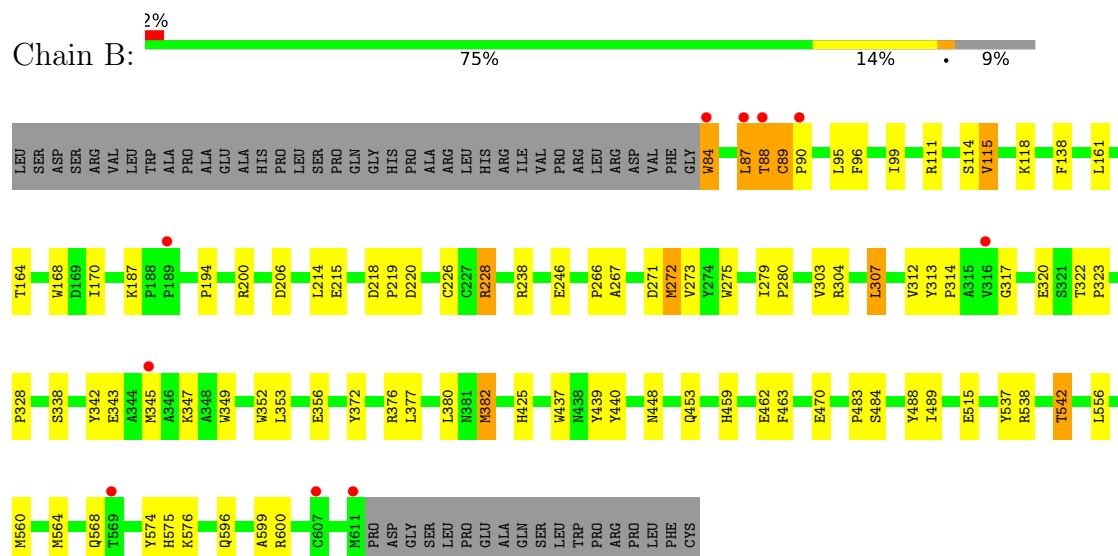
3 Residue-property plots [i](#)

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

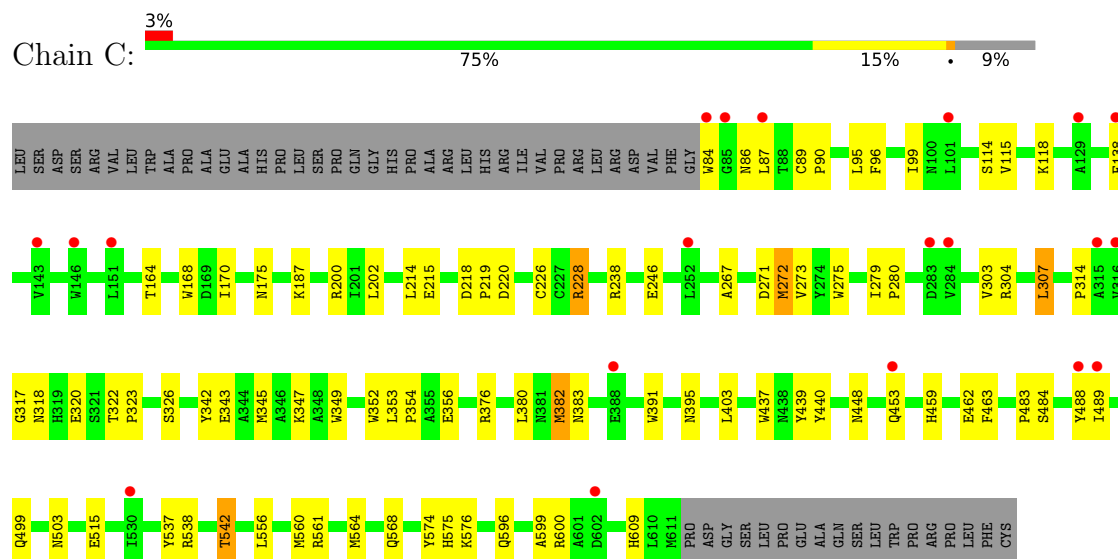
• Molecule 1: Spingomyelin phosphodiesterase



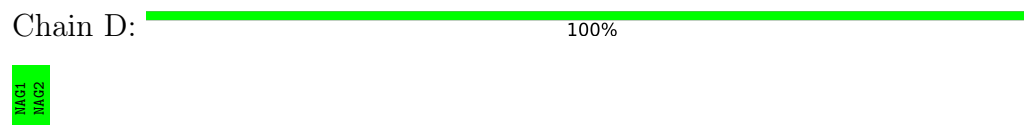
• Molecule 1: Spingomyelin phosphodiesterase



- Molecule 1: Sphingomyelin phosphodiesterase



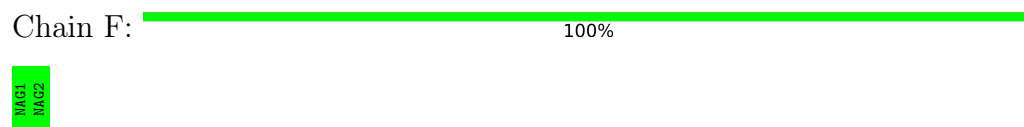
- Molecule 2: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose



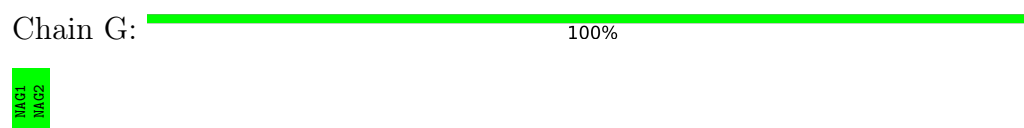
- Molecule 2: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose



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



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



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

Chain I:  100%

MAG1
MAG2

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

Chain H:  33% 33% 33%

MAG1
MAG2
BMA3

4 Data and refinement statistics

Property	Value	Source
Space group	I 2 2 2	Depositor
Cell constants a, b, c, α , β , γ	191.02Å 230.87Å 252.32Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	44.13 – 3.65 44.13 – 3.65	Depositor EDS
% Data completeness (in resolution range)	98.6 (44.13-3.65) 98.8 (44.13-3.65)	Depositor EDS
R_{merge}	0.23	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.34 (at 3.66Å)	Xtriage
Refinement program	PHENIX (dev_2229: ???)	Depositor
R, R_{free}	0.248 , 0.254 0.252 , 0.262	Depositor DCC
R_{free} test set	2000 reflections (3.25%)	wwPDB-VP
Wilson B-factor (Å ²)	188.7	Xtriage
Anisotropy	0.283	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 220.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.38$, $\langle L^2 \rangle = 0.22$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	12883	wwPDB-VP
Average B, all atoms (Å ²)	237.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.54% 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: ZN, NAG, MAN, BMA

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	0.22	0/4307	0.56	0/5904
1	B	0.20	0/4307	0.57	1/5904 (0.0%)
1	C	0.18	0/4307	0.51	0/5904
All	All	0.20	0/12921	0.55	1/17712 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	84	TRP	N-CA-C	5.50	126.39	111.00

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	4161	0	4027	80	0
1	B	4161	0	4024	59	0
1	C	4161	0	4032	61	0
2	D	28	0	25	0	0
2	E	28	0	25	2	0
2	F	28	0	25	0	0
2	G	28	0	25	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	I	28	0	25	0	0
3	H	39	0	34	3	0
4	A	70	0	63	9	0
4	B	56	0	51	2	0
4	C	56	0	52	10	0
5	A	11	0	10	2	0
5	B	11	0	10	2	0
6	A	11	0	10	0	0
7	A	2	0	0	0	0
7	B	2	0	0	0	0
7	C	2	0	0	0	0
All	All	12883	0	12438	208	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 8.

All (208) 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:86:ASN:HD21	4:A:701:NAG:C1	1.07	1.64
1:C:86:ASN:HD21	4:C:701:NAG:C1	1.24	1.47
1:C:395:ASN:HD21	4:C:703:NAG:C1	1.27	1.45
4:A:709:NAG:O4	4:A:710:NAG:C1	1.63	1.45
1:A:86:ASN:ND2	4:A:701:NAG:C1	1.88	1.33
5:A:707:BMA:C1	2:E:2:NAG:O4	1.81	1.29
1:C:175:ASN:ND2	4:C:702:NAG:C1	1.99	1.25
1:C:503:ASN:HD21	4:C:704:NAG:C1	1.51	1.23
4:B:708:NAG:O4	5:B:709:BMA:C1	1.91	1.19
1:C:86:ASN:ND2	4:C:701:NAG:C1	2.09	1.14
1:C:395:ASN:ND2	4:C:703:NAG:C1	2.10	1.13
1:A:86:ASN:ND2	4:A:701:NAG:O5	1.82	1.10
3:H:2:NAG:O3	3:H:3:BMA:H2	1.54	1.05
1:C:175:ASN:HD21	4:C:702:NAG:C1	1.64	1.02
1:C:503:ASN:ND2	4:C:704:NAG:C1	2.22	1.01
4:A:709:NAG:C4	4:A:710:NAG:C1	2.45	0.95
1:C:175:ASN:HD22	4:C:702:NAG:C1	1.82	0.87
4:A:709:NAG:O4	4:A:710:NAG:C2	2.25	0.84
4:B:708:NAG:HO4	5:B:709:BMA:C1	1.88	0.83
3:H:2:NAG:H4	3:H:3:BMA:O2	1.81	0.79
3:H:2:NAG:O3	3:H:3:BMA:C2	2.27	0.79
1:A:86:ASN:CG	4:A:701:NAG:C1	2.60	0.75

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:218:ASP:OD1	1:C:238:ARG:NH2	2.21	0.74
1:C:515:GLU:HG2	1:C:538:ARG:HG2	1.71	0.73
1:C:462:GLU:HB2	1:C:596:GLN:HE22	1.52	0.73
1:C:537:TYR:HB2	1:C:542:THR:HG21	1.69	0.72
1:B:342:TYR:HA	1:B:345:MET:HE2	1.71	0.72
1:B:87:LEU:O	1:B:88:THR:OG1	2.09	0.71
1:B:515:GLU:HG2	1:B:538:ARG:HG2	1.73	0.71
1:B:380:LEU:HB3	1:B:382:MET:HE3	1.73	0.70
1:C:86:ASN:ND2	4:C:701:NAG:O5	2.21	0.70
1:C:187:LYS:H	1:C:448:ASN:HD21	1.39	0.70
1:A:342:TYR:HA	1:A:345:MET:HE2	1.75	0.69
1:C:342:TYR:HA	1:C:345:MET:HE2	1.74	0.69
1:B:200:ARG:NH1	1:B:267:ALA:O	2.24	0.68
1:B:273:VAL:HG21	1:B:307:LEU:HD22	1.75	0.68
1:A:380:LEU:HB3	1:A:382:MET:HE3	1.76	0.67
1:B:382:MET:HA	1:B:382:MET:HE2	1.77	0.67
1:B:462:GLU:HB2	1:B:596:GLN:HE22	1.60	0.67
1:C:95:LEU:O	1:C:99:ILE:HG12	1.95	0.66
1:A:463:PHE:H	1:A:596:GLN:HE22	1.42	0.66
1:A:382:MET:HE2	1:A:382:MET:HA	1.78	0.66
1:A:564:MET:HE3	1:A:568:GLN:HE21	1.61	0.65
1:B:537:TYR:HB2	1:B:542:THR:HG21	1.79	0.64
1:A:187:LYS:H	1:A:448:ASN:HD21	1.44	0.63
1:C:382:MET:HE2	1:C:382:MET:HA	1.79	0.63
1:C:564:MET:HE3	1:C:568:GLN:HE21	1.64	0.63
1:A:515:GLU:HG2	1:A:538:ARG:HG2	1.80	0.63
1:B:89:CYS:HB3	1:B:90:PRO:HD3	1.81	0.63
1:B:439:TYR:OH	1:B:453:GLN:NE2	2.33	0.62
1:A:537:TYR:HB2	1:A:542:THR:HG21	1.81	0.62
1:A:138:PHE:CE1	1:A:323:PRO:HB3	2.34	0.61
1:A:218:ASP:OD1	1:A:238:ARG:NH2	2.30	0.61
4:A:709:NAG:H4	4:A:710:NAG:C1	2.28	0.61
1:C:273:VAL:HG21	1:C:307:LEU:HD22	1.82	0.61
1:B:440:TYR:HE1	1:B:599:ALA:HB3	1.65	0.60
1:C:380:LEU:HB3	1:C:382:MET:HE3	1.83	0.60
1:B:564:MET:HE3	1:B:568:GLN:HE21	1.67	0.60
1:A:273:VAL:HG21	1:A:307:LEU:HD22	1.85	0.59
1:B:95:LEU:O	1:B:99:ILE:HG12	2.02	0.59
1:C:114:SER:O	1:C:118:LYS:HG2	2.01	0.59
1:A:538:ARG:O	1:A:542:THR:HG23	2.02	0.59
1:A:95:LEU:O	1:A:99:ILE:HG12	2.03	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:318:ASN:HA	1:C:383:ASN:HD21	1.68	0.58
1:A:114:SER:O	1:A:118:LYS:HG2	2.03	0.58
1:A:207:LEU:HA	1:A:485:ALA:HB3	1.85	0.58
1:B:304:ARG:HG3	1:B:352:TRP:CZ2	2.39	0.58
1:C:499:GLN:NE2	1:C:515:GLU:OE2	2.35	0.58
1:A:200:ARG:NH1	1:A:267:ALA:O	2.36	0.57
1:B:463:PHE:H	1:B:596:GLN:HE22	1.52	0.57
1:C:463:PHE:H	1:C:596:GLN:HE22	1.51	0.57
1:C:138:PHE:CE1	1:C:323:PRO:HB3	2.40	0.56
1:A:271:ASP:C	1:A:272:MET:HE3	2.29	0.56
1:B:114:SER:O	1:B:118:LYS:HG2	2.05	0.56
1:B:271:ASP:C	1:B:272:MET:HE3	2.30	0.56
1:B:538:ARG:O	1:B:542:THR:HG23	2.06	0.56
5:A:707:BMA:C1	2:E:2:NAG:C4	2.83	0.55
1:B:218:ASP:OD1	1:B:238:ARG:NH2	2.35	0.55
1:C:202:LEU:HB3	1:C:273:VAL:HG22	1.88	0.55
1:B:138:PHE:CE1	1:B:323:PRO:HB3	2.42	0.54
1:A:462:GLU:HB2	1:A:596:GLN:HE22	1.72	0.54
1:A:252:LEU:HD11	1:A:485:ALA:O	2.07	0.54
1:C:538:ARG:O	1:C:542:THR:HG23	2.07	0.54
1:B:187:LYS:H	1:B:448:ASN:HD21	1.56	0.54
1:C:556:LEU:HG	1:C:560:MET:HE2	1.90	0.53
1:C:168:TRP:CH2	1:C:170:ILE:HD12	2.43	0.53
1:A:279:ILE:HB	1:A:280:PRO:HD3	1.90	0.53
1:A:440:TYR:CD1	1:A:600:ARG:HB2	2.43	0.53
1:C:343:GLU:O	1:C:347:LYS:HD3	2.09	0.53
1:C:561:ARG:NH1	1:C:609:HIS:O	2.42	0.53
1:A:220:ASP:HB3	1:A:238:ARG:NH2	2.24	0.52
1:C:271:ASP:HB3	1:C:272:MET:HE3	1.90	0.52
1:B:88:THR:O	1:B:161:LEU:HD13	2.09	0.52
1:B:96:PHE:HA	1:B:99:ILE:HB	1.92	0.52
1:B:343:GLU:O	1:B:347:LYS:HD3	2.10	0.52
1:C:440:TYR:HE1	1:C:599:ALA:HB3	1.75	0.52
1:C:219:PRO:HB3	1:C:246:GLU:HG2	1.92	0.52
1:A:96:PHE:HA	1:A:99:ILE:HB	1.92	0.51
1:B:220:ASP:HB3	1:B:238:ARG:NH2	2.25	0.51
1:A:168:TRP:CH2	1:A:170:ILE:HD12	2.46	0.51
1:A:318:ASN:HA	1:A:383:ASN:HD21	1.75	0.51
1:A:483:PRO:HB2	1:A:576:LYS:HD3	1.93	0.51
1:B:275:TRP:O	1:B:314:PRO:HA	2.11	0.51
1:C:483:PRO:HB2	1:C:576:LYS:HD3	1.93	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:463:PHE:H	1:A:596:GLN:NE2	2.09	0.50
1:A:89:CYS:HB3	1:A:90:PRO:HD3	1.93	0.50
1:A:372:TYR:CE1	1:A:503:ASN:HB2	2.46	0.50
1:B:312:VAL:HB	1:B:352:TRP:CD2	2.46	0.49
1:B:437:TRP:CD2	1:B:600:ARG:HD2	2.47	0.49
1:B:556:LEU:HG	1:B:560:MET:HE2	1.94	0.49
1:A:314:PRO:HG2	1:A:349:TRP:CD2	2.48	0.49
1:A:343:GLU:O	1:A:347:LYS:HD3	2.13	0.49
1:C:314:PRO:HG2	1:C:349:TRP:CD2	2.48	0.49
1:A:159:LEU:HB2	1:A:168:TRP:CE3	2.47	0.49
1:C:275:TRP:O	1:C:314:PRO:HA	2.13	0.48
1:B:322:THR:HA	1:B:323:PRO:C	2.38	0.48
1:C:271:ASP:C	1:C:272:MET:HE3	2.38	0.48
1:A:304:ARG:HG3	1:A:352:TRP:CZ2	2.49	0.48
1:B:194:PRO:HG3	1:C:538:ARG:HD3	1.95	0.47
1:B:380:LEU:CB	1:B:382:MET:HE3	2.43	0.47
1:C:304:ARG:HG3	1:C:352:TRP:CZ2	2.49	0.47
1:C:96:PHE:HA	1:C:99:ILE:HB	1.95	0.47
1:C:463:PHE:H	1:C:596:GLN:NE2	2.11	0.47
1:A:275:TRP:O	1:A:314:PRO:HA	2.14	0.47
1:A:459:HIS:HA	1:A:484:SER:HB3	1.96	0.47
1:A:219:PRO:HB3	1:A:246:GLU:HG2	1.95	0.47
1:C:220:ASP:HB3	1:C:238:ARG:NH2	2.30	0.47
1:A:439:TYR:OH	1:A:453:GLN:NE2	2.48	0.47
1:B:314:PRO:HG2	1:B:349:TRP:CD2	2.49	0.47
1:A:387:ARG:HH22	1:A:601:ALA:HB1	1.80	0.47
1:A:428:PRO:HD3	1:A:455:PHE:CZ	2.50	0.47
1:C:488:TYR:HA	1:C:489:ILE:HA	1.61	0.47
1:B:515:GLU:CG	1:B:538:ARG:HG2	2.44	0.47
1:C:322:THR:HA	1:C:323:PRO:C	2.40	0.47
1:A:194:PRO:HG3	1:B:538:ARG:HD3	1.96	0.46
1:C:380:LEU:CB	1:C:382:MET:HE3	2.45	0.46
1:C:439:TYR:OH	1:C:453:GLN:NE2	2.47	0.46
1:C:89:CYS:HB3	1:C:90:PRO:HD3	1.96	0.46
1:C:200:ARG:NH1	1:C:267:ALA:O	2.40	0.46
1:A:380:LEU:CB	1:A:382:MET:HE3	2.44	0.46
1:B:219:PRO:HB3	1:B:246:GLU:HG2	1.98	0.46
1:C:279:ILE:HB	1:C:280:PRO:HD3	1.97	0.46
1:A:437:TRP:CD2	1:A:600:ARG:HD2	2.51	0.46
1:C:214:LEU:HD23	1:C:214:LEU:HA	1.78	0.46
1:B:272:MET:HG2	1:B:313:TYR:CE2	2.50	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:214:LEU:HD23	1:A:214:LEU:HA	1.67	0.45
1:B:214:LEU:HD23	1:B:214:LEU:HA	1.83	0.45
1:A:86:ASN:OD1	4:A:701:NAG:C1	2.65	0.45
1:A:312:VAL:HB	1:A:352:TRP:CD2	2.52	0.45
1:A:561:ARG:HD3	1:A:610:LEU:HD23	1.97	0.45
1:B:483:PRO:HB2	1:B:576:LYS:HD3	1.98	0.45
1:A:261:LEU:HD11	1:A:303:VAL:HG12	1.98	0.44
1:A:271:ASP:HB3	1:A:272:MET:HE3	1.99	0.44
1:A:466:PHE:HB2	1:A:477:ALA:HB3	1.99	0.44
1:A:308:GLY:HA3	1:A:309:PRO:HD3	1.78	0.44
1:A:191:PRO:HD3	1:B:266:PRO:HB3	1.98	0.44
1:A:244:TRP:CZ2	1:A:531:PRO:HG3	2.53	0.44
1:A:322:THR:HA	1:A:323:PRO:C	2.41	0.44
1:A:428:PRO:HD3	1:A:455:PHE:CE2	2.51	0.44
1:B:488:TYR:HA	1:B:489:ILE:HA	1.63	0.44
1:A:85:GLY:O	1:A:88:THR:HG23	2.18	0.44
1:A:201:ILE:HD12	1:A:272:MET:HB2	1.99	0.44
1:B:463:PHE:H	1:B:596:GLN:NE2	2.14	0.43
1:C:459:HIS:HA	1:C:484:SER:HB3	2.00	0.43
1:A:488:TYR:HA	1:A:489:ILE:HA	1.60	0.43
1:B:313:TYR:HD2	1:B:377:LEU:HD13	1.84	0.43
1:B:440:TYR:CD1	1:B:600:ARG:HB2	2.53	0.43
1:A:87:LEU:HB3	1:A:91:ILE:HD12	2.01	0.43
1:B:574:TYR:HD1	1:B:575:HIS:CE1	2.36	0.43
1:A:207:LEU:HB3	1:A:257:LEU:HD11	2.01	0.43
1:B:380:LEU:HB3	1:B:382:MET:CE	2.47	0.43
1:A:244:TRP:CE2	1:A:531:PRO:HG3	2.54	0.43
1:A:556:LEU:HG	1:A:560:MET:HE2	2.00	0.43
1:A:361:LEU:HD12	1:A:365:GLY:HA2	2.02	0.42
1:A:387:ARG:NH2	1:A:601:ALA:HB1	2.34	0.42
1:B:317:GLY:O	1:B:320:GLU:HG2	2.20	0.42
1:A:318:ASN:CA	1:A:383:ASN:HD21	2.33	0.42
1:C:574:TYR:HD1	1:C:575:HIS:CE1	2.37	0.42
1:A:328:PRO:HD2	1:A:338:SER:OG	2.20	0.42
1:B:470:GLU:OE1	1:B:470:GLU:N	2.46	0.42
1:A:294:ARG:HG3	1:A:295:ALA:N	2.34	0.42
1:A:320:GLU:HB2	1:A:341:LEU:HD21	2.01	0.42
1:A:328:PRO:HA	1:A:329:PRO:HD3	1.91	0.42
1:B:206:ASP:OD1	1:B:425:HIS:HE1	2.03	0.42
1:C:403:LEU:HD23	1:C:403:LEU:HA	1.94	0.42
1:B:459:HIS:HA	1:B:484:SER:HB3	2.02	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:168:TRP:CH2	1:B:170:ILE:HD12	2.55	0.41
1:A:203:PHE:CE1	1:A:454:PHE:HB3	2.55	0.41
1:B:272:MET:HE1	1:B:372:TYR:CZ	2.55	0.41
1:B:279:ILE:HB	1:B:280:PRO:HD3	2.03	0.41
1:C:226:CYS:O	1:C:228:ARG:HG2	2.20	0.41
1:A:247:TYR:CZ	1:A:521:LEU:HB2	2.54	0.41
1:C:326:SER:HG	1:C:391:TRP:HE1	1.65	0.41
1:C:440:TYR:CD1	1:C:600:ARG:HB2	2.56	0.41
1:A:470:GLU:OE1	1:A:470:GLU:N	2.45	0.41
1:B:226:CYS:O	1:B:228:ARG:HG2	2.20	0.41
1:A:96:PHE:HA	1:A:99:ILE:CG1	2.50	0.41
1:B:328:PRO:HD2	1:B:338:SER:OG	2.21	0.41
1:A:437:TRP:CG	1:A:600:ARG:HD2	2.56	0.41
1:B:271:ASP:HB3	1:B:272:MET:HE3	2.03	0.41
1:C:317:GLY:O	1:C:320:GLU:HG2	2.21	0.41
1:C:437:TRP:CG	1:C:600:ARG:HD2	2.56	0.41
1:A:251:ASP:OD2	1:A:282:HIS:HD2	2.05	0.40
1:A:499:GLN:HE21	1:A:513:ASP:HB3	1.86	0.40
1:C:271:ASP:CB	1:C:272:MET:HE3	2.51	0.40
1:B:111:ARG:O	1:B:115:VAL:HG12	2.21	0.40
1:A:293:LEU:HD23	1:A:293:LEU:HA	1.88	0.40
1:B:437:TRP:CG	1:B:600:ARG:HD2	2.56	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	526/583 (90%)	503 (96%)	22 (4%)	1 (0%)	43 71
1	B	526/583 (90%)	502 (95%)	22 (4%)	2 (0%)	30 59
1	C	526/583 (90%)	503 (96%)	22 (4%)	1 (0%)	43 71

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
All	All	1578/1749 (90%)	1508 (96%)	66 (4%)	4 (0%)	36 64

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	88	THR
1	B	89	CYS
1	A	354	PRO
1	C	354	PRO

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	448/495 (90%)	434 (97%)	14 (3%)	35 55
1	B	448/495 (90%)	434 (97%)	14 (3%)	35 55
1	C	448/495 (90%)	434 (97%)	14 (3%)	35 55
All	All	1344/1485 (90%)	1302 (97%)	42 (3%)	35 55

All (42) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	84	TRP
1	A	87	LEU
1	A	115	VAL
1	A	164	THR
1	A	215	GLU
1	A	228	ARG
1	A	272	MET
1	A	303	VAL
1	A	307	LEU
1	A	353	LEU
1	A	356	GLU
1	A	376	ARG

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Mol	Chain	Res	Type
1	A	382	MET
1	A	542	THR
1	B	84	TRP
1	B	87	LEU
1	B	115	VAL
1	B	164	THR
1	B	215	GLU
1	B	228	ARG
1	B	272	MET
1	B	303	VAL
1	B	307	LEU
1	B	353	LEU
1	B	356	GLU
1	B	376	ARG
1	B	382	MET
1	B	542	THR
1	C	84	TRP
1	C	87	LEU
1	C	115	VAL
1	C	164	THR
1	C	215	GLU
1	C	228	ARG
1	C	272	MET
1	C	303	VAL
1	C	307	LEU
1	C	353	LEU
1	C	356	GLU
1	C	376	ARG
1	C	382	MET
1	C	542	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (45) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	86	ASN
1	A	208	HIS
1	A	282	HIS
1	A	287	GLN
1	A	336	HIS
1	A	383	ASN
1	A	411	GLN
1	A	430	HIS

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Mol	Chain	Res	Type
1	A	448	ASN
1	A	453	GLN
1	A	499	GLN
1	A	514	HIS
1	A	568	GLN
1	A	578	HIS
1	A	596	GLN
1	B	287	GLN
1	B	383	ASN
1	B	411	GLN
1	B	430	HIS
1	B	448	ASN
1	B	453	GLN
1	B	514	HIS
1	B	568	GLN
1	B	575	HIS
1	B	578	HIS
1	B	596	GLN
1	C	86	ASN
1	C	136	HIS
1	C	175	ASN
1	C	208	HIS
1	C	287	GLN
1	C	336	HIS
1	C	383	ASN
1	C	395	ASN
1	C	411	GLN
1	C	448	ASN
1	C	453	GLN
1	C	499	GLN
1	C	503	ASN
1	C	509	HIS
1	C	514	HIS
1	C	568	GLN
1	C	575	HIS
1	C	578	HIS
1	C	596	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 ⓘ

13 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	D	1	1,2	14,14,15	0.20	0	17,19,21	0.41	0
2	NAG	D	2	2	14,14,15	0.31	0	17,19,21	0.59	0
2	NAG	E	1	1,2	14,14,15	0.15	0	17,19,21	0.48	0
2	NAG	E	2	2	14,14,15	0.43	0	17,19,21	0.63	0
2	NAG	F	1	1,2	14,14,15	0.22	0	17,19,21	0.43	0
2	NAG	F	2	2	14,14,15	0.22	0	17,19,21	0.52	0
2	NAG	G	1	1,2	14,14,15	0.18	0	17,19,21	0.44	0
2	NAG	G	2	2	14,14,15	0.22	0	17,19,21	0.43	0
3	NAG	H	1	1,3	14,14,15	0.28	0	17,19,21	0.47	0
3	NAG	H	2	3	14,14,15	0.56	0	17,19,21	1.21	2 (11%)
3	BMA	H	3	3	11,11,12	0.26	0	15,15,17	0.52	0
2	NAG	I	1	1,2	14,14,15	0.26	0	17,19,21	0.39	0
2	NAG	I	2	2	14,14,15	0.31	0	17,19,21	0.38	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	D	1	1,2	-	2/6/23/26	0/1/1/1
2	NAG	D	2	2	-	2/6/23/26	0/1/1/1
2	NAG	E	1	1,2	-	0/6/23/26	0/1/1/1
2	NAG	E	2	2	-	1/6/23/26	0/1/1/1
2	NAG	F	1	1,2	-	1/6/23/26	0/1/1/1
2	NAG	F	2	2	-	2/6/23/26	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	NAG	G	1	1,2	-	0/6/23/26	0/1/1/1
2	NAG	G	2	2	-	0/6/23/26	0/1/1/1
3	NAG	H	1	1,3	-	0/6/23/26	0/1/1/1
3	NAG	H	2	3	-	5/6/23/26	0/1/1/1
3	BMA	H	3	3	-	0/2/19/22	0/1/1/1
2	NAG	I	1	1,2	-	0/6/23/26	0/1/1/1
2	NAG	I	2	2	-	0/6/23/26	0/1/1/1

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	H	2	NAG	O5-C1-C2	-2.88	106.83	111.29
3	H	2	NAG	C2-N2-C7	-2.09	120.10	122.90

There are no chirality outliers.

All (13) torsion outliers are listed below:

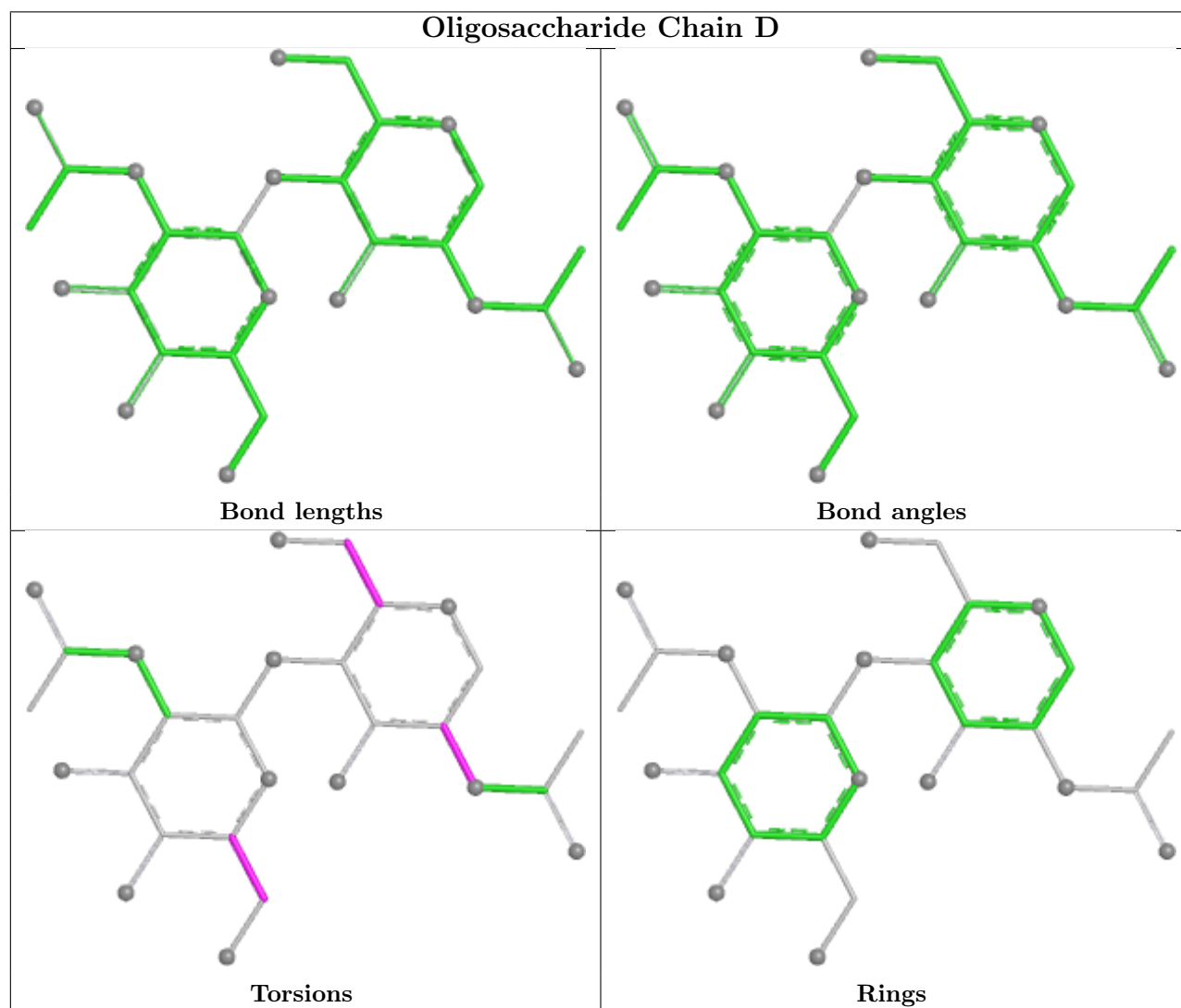
Mol	Chain	Res	Type	Atoms
3	H	2	NAG	C8-C7-N2-C2
3	H	2	NAG	O7-C7-N2-C2
2	D	2	NAG	C4-C5-C6-O6
2	F	2	NAG	C4-C5-C6-O6
2	D	2	NAG	O5-C5-C6-O6
2	F	2	NAG	O5-C5-C6-O6
3	H	2	NAG	O5-C5-C6-O6
3	H	2	NAG	C4-C5-C6-O6
3	H	2	NAG	C3-C2-N2-C7
2	D	1	NAG	C1-C2-N2-C7
2	E	2	NAG	C1-C2-N2-C7
2	F	1	NAG	C4-C5-C6-O6
2	D	1	NAG	C4-C5-C6-O6

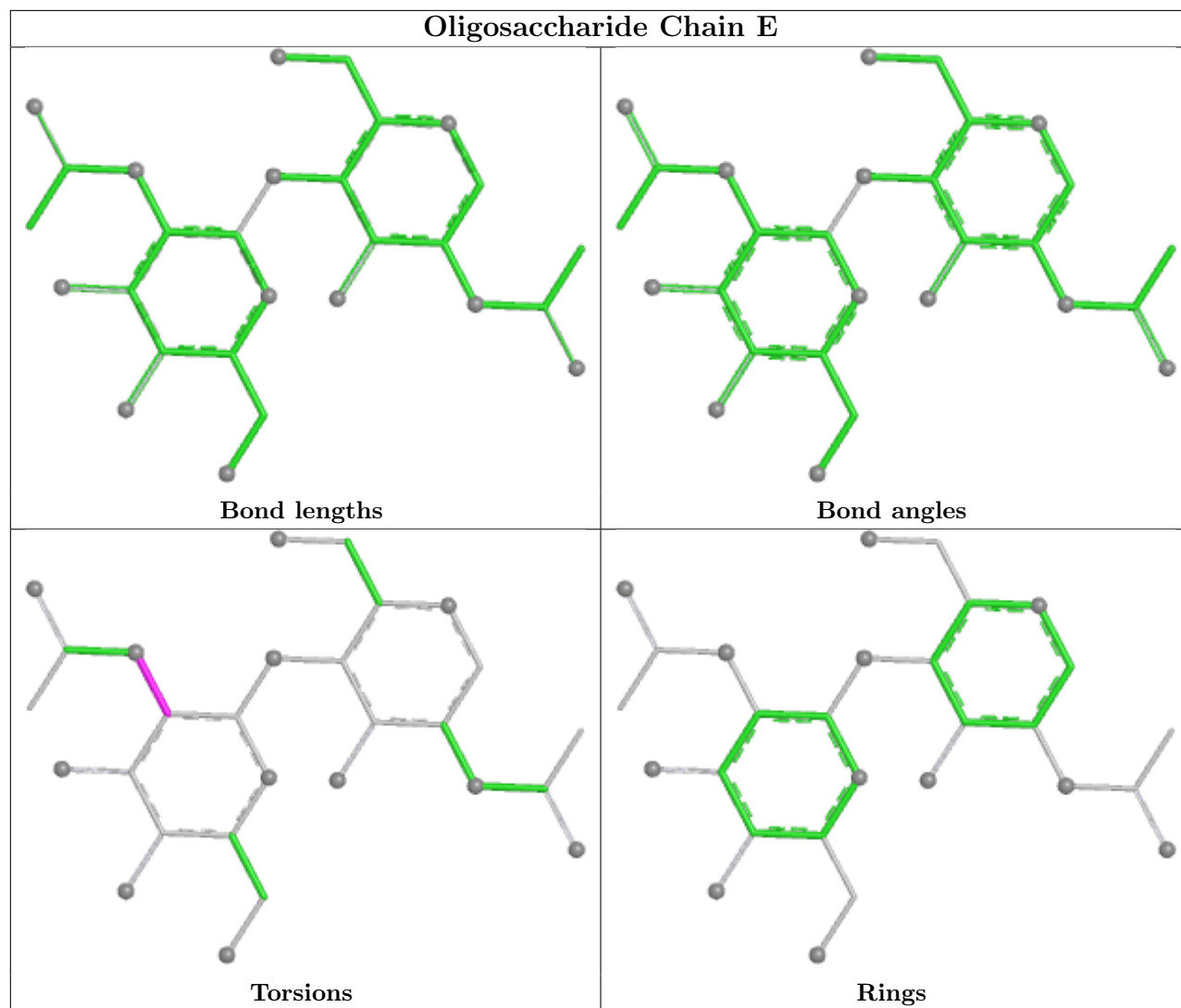
There are no ring outliers.

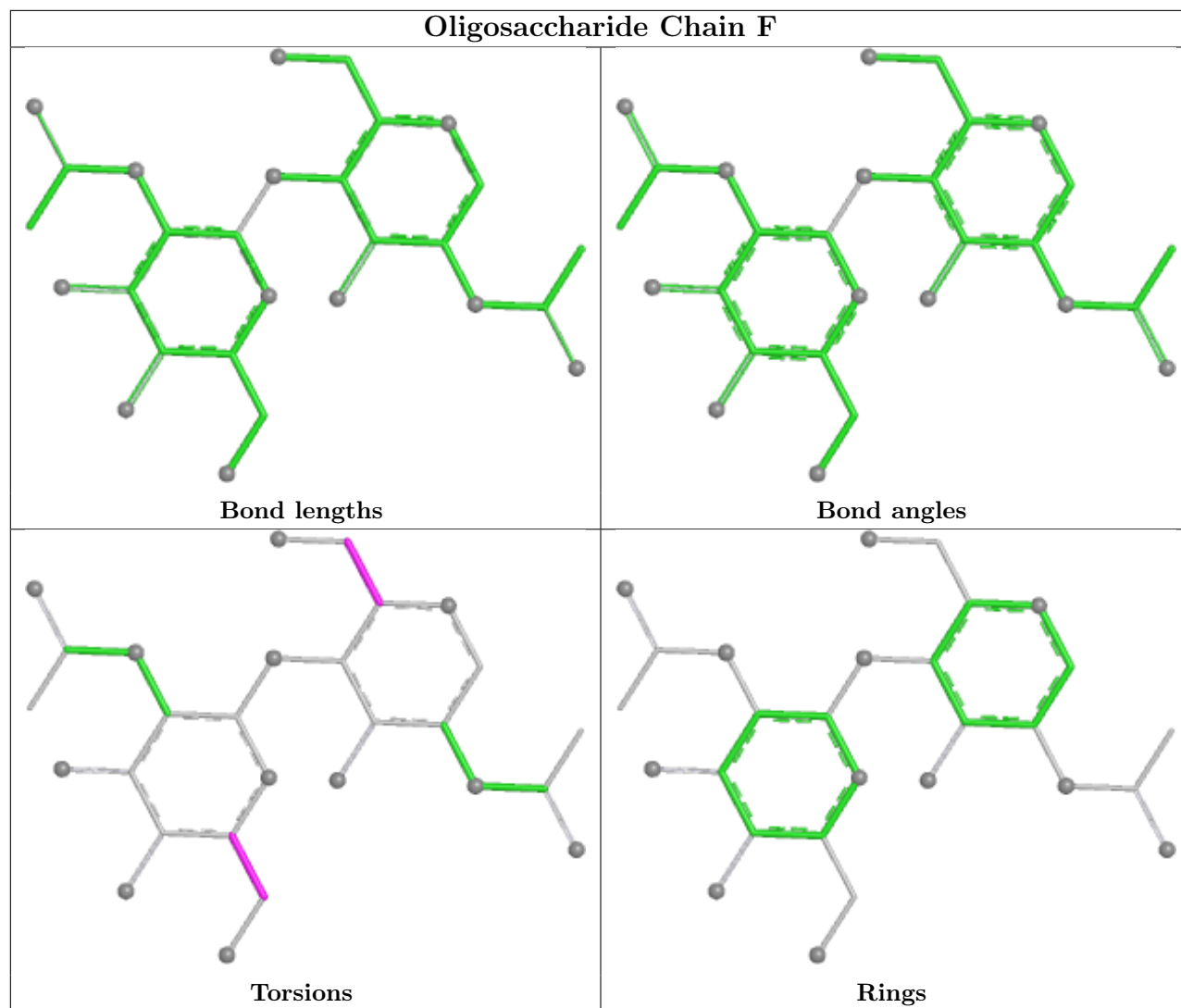
3 monomers are involved in 5 short contacts:

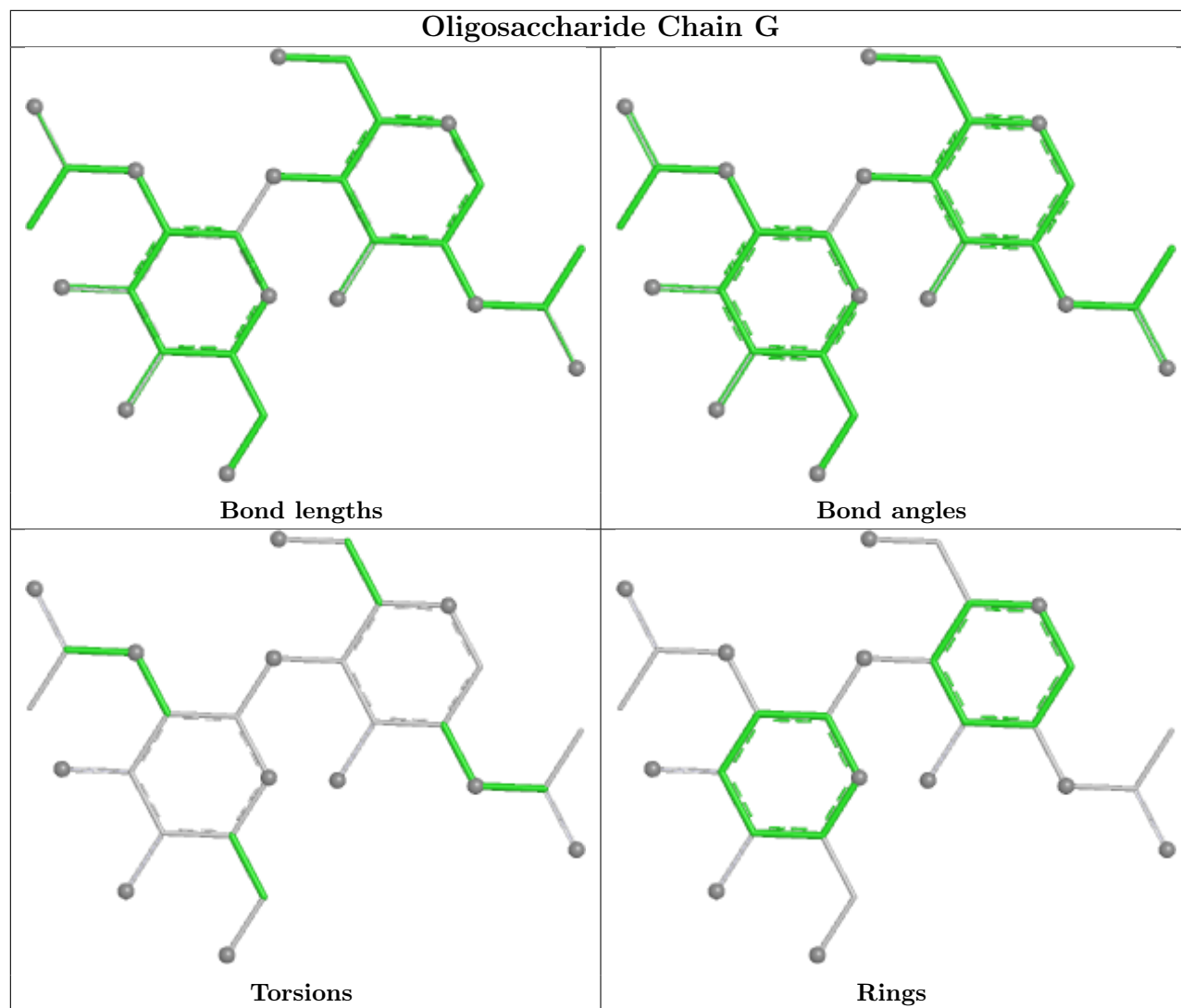
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	H	3	BMA	3	0
2	E	2	NAG	2	0
3	H	2	NAG	3	0

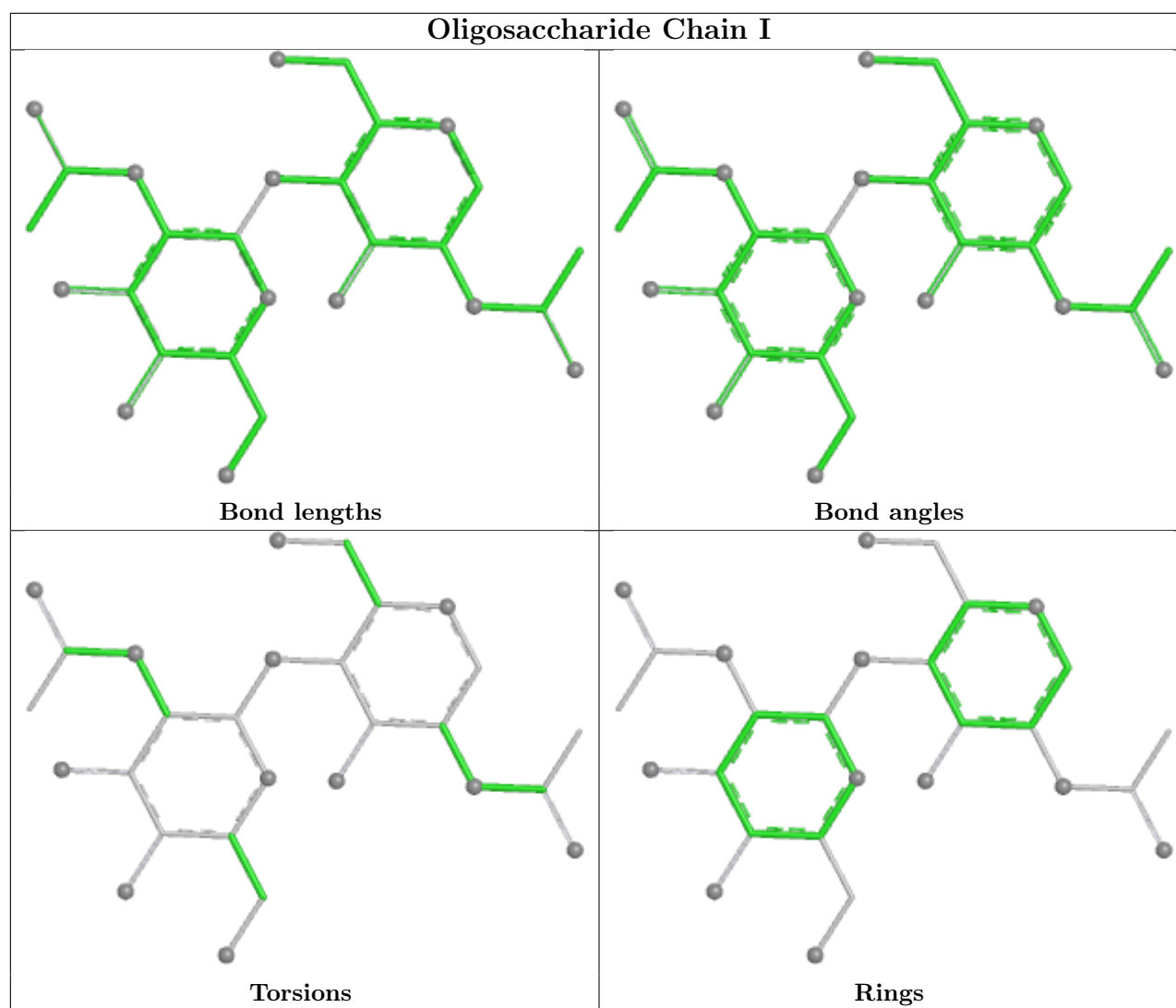
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for oligosaccharide.

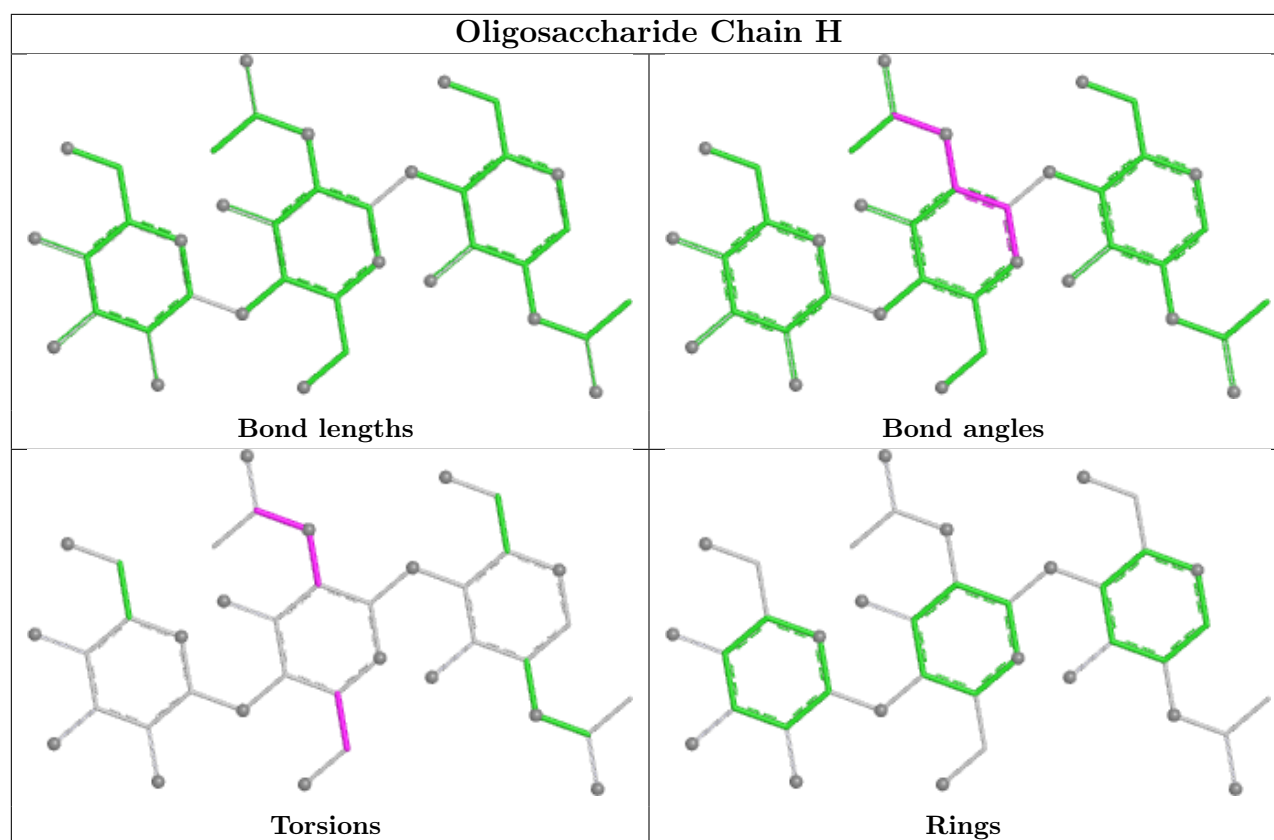












5.6 Ligand geometry [i](#)

Of 22 ligands modelled in this entry, 6 are monoatomic - leaving 16 for Mogul analysis.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
4	NAG	A	709	1	14,14,15	0.25	0	17,19,21	0.52	0
4	NAG	C	702	-	14,14,15	0.30	0	17,19,21	0.64	0
4	NAG	B	708	-	14,14,15	0.40	0	17,19,21	0.52	0
4	NAG	C	701	-	14,14,15	0.94	1 (7%)	17,19,21	0.57	0
4	NAG	B	707	1	14,14,15	0.25	0	17,19,21	0.47	0
4	NAG	A	708	-	14,14,15	0.39	0	17,19,21	0.53	0
4	NAG	B	701	1	14,14,15	0.22	0	17,19,21	0.48	0
5	BMA	B	709	-	11,11,12	1.10	1 (9%)	15,15,17	1.40	3 (20%)
6	MAN	A	711	-	11,11,12	0.83	0	15,15,17	0.94	1 (6%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
4	NAG	A	702	1	14,14,15	0.24	0	17,19,21	0.55	0
4	NAG	A	710	-	14,14,15	0.38	0	17,19,21	0.40	0
4	NAG	C	704	-	14,14,15	0.77	1 (7%)	17,19,21	0.61	0
4	NAG	A	701	-	14,14,15	0.66	1 (7%)	17,19,21	0.46	0
4	NAG	B	702	-	14,14,15	0.24	0	17,19,21	0.54	0
4	NAG	C	703	-	14,14,15	0.68	1 (7%)	17,19,21	0.66	1 (5%)
5	BMA	A	707	-	11,11,12	1.14	2 (18%)	15,15,17	1.39	3 (20%)

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	NAG	A	709	1	-	0/6/23/26	0/1/1/1
4	NAG	C	702	-	-	2/6/23/26	0/1/1/1
4	NAG	B	708	-	-	0/6/23/26	0/1/1/1
4	NAG	C	701	-	-	0/6/23/26	0/1/1/1
4	NAG	B	707	1	-	2/6/23/26	0/1/1/1
4	NAG	A	708	-	-	2/6/23/26	0/1/1/1
4	NAG	B	701	1	-	0/6/23/26	0/1/1/1
5	BMA	B	709	-	-	0/2/19/22	0/1/1/1
6	MAN	A	711	-	-	2/2/19/22	0/1/1/1
4	NAG	A	702	1	-	2/6/23/26	0/1/1/1
4	NAG	A	710	-	-	0/6/23/26	0/1/1/1
4	NAG	C	704	-	-	2/6/23/26	0/1/1/1
4	NAG	A	701	-	-	0/6/23/26	0/1/1/1
4	NAG	B	702	-	-	2/6/23/26	0/1/1/1
4	NAG	C	703	-	-	3/6/23/26	0/1/1/1
5	BMA	A	707	-	-	0/2/19/22	0/1/1/1

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	C	701	NAG	C1-C2	3.16	1.56	1.52
5	B	709	BMA	C1-C2	2.49	1.58	1.52
5	A	707	BMA	C2-C3	2.42	1.56	1.52
4	C	704	NAG	C1-C2	2.41	1.55	1.52
5	A	707	BMA	C1-C2	2.34	1.57	1.52

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	701	NAG	C1-C2	2.33	1.55	1.52
4	C	703	NAG	O5-C1	2.13	1.47	1.43

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	B	709	BMA	C1-O5-C5	3.50	116.88	112.19
5	A	707	BMA	C1-O5-C5	2.97	116.16	112.19
5	A	707	BMA	C1-C2-C3	2.95	113.94	109.64
5	B	709	BMA	C1-C2-C3	2.53	113.33	109.64
5	B	709	BMA	O2-C2-C3	-2.50	104.97	110.15
6	A	711	MAN	C1-O5-C5	2.28	115.25	112.19
4	C	703	NAG	C1-O5-C5	2.24	115.18	112.19
5	A	707	BMA	O2-C2-C3	-2.13	105.73	110.15

There are no chirality outliers.

All (17) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	B	702	NAG	O5-C5-C6-O6
4	B	707	NAG	O5-C5-C6-O6
4	C	702	NAG	C4-C5-C6-O6
4	B	702	NAG	C4-C5-C6-O6
4	C	702	NAG	O5-C5-C6-O6
4	B	707	NAG	C4-C5-C6-O6
4	C	703	NAG	O5-C5-C6-O6
4	A	702	NAG	C4-C5-C6-O6
4	A	702	NAG	O5-C5-C6-O6
6	A	711	MAN	O5-C5-C6-O6
4	C	703	NAG	C4-C5-C6-O6
4	A	708	NAG	C4-C5-C6-O6
4	A	708	NAG	O5-C5-C6-O6
4	C	704	NAG	C4-C5-C6-O6
4	C	704	NAG	O5-C5-C6-O6
4	C	703	NAG	C1-C2-N2-C7
6	A	711	MAN	C4-C5-C6-O6

There are no ring outliers.

10 monomers are involved in 23 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	709	NAG	4	0
4	C	702	NAG	3	0
4	B	708	NAG	2	0
4	C	701	NAG	3	0
5	B	709	BMA	2	0
4	A	710	NAG	4	0
4	C	704	NAG	2	0
4	A	701	NAG	5	0
4	C	703	NAG	2	0
5	A	707	BMA	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

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	528/583 (90%)	-0.15	4 (0%) 82 60	147, 194, 235, 267	0
1	B	528/583 (90%)	-0.05	10 (1%) 66 40	167, 211, 266, 321	0
1	C	528/583 (90%)	0.09	20 (3%) 44 26	215, 302, 350, 413	0
All	All	1584/1749 (90%)	-0.04	34 (2%) 63 38	147, 221, 335, 413	0

All (34) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	489	ILE	4.2
1	B	611	MET	3.8
1	B	84	TRP	3.7
1	A	84	TRP	3.1
1	B	345	MET	3.1
1	B	607	CYS	2.9
1	C	488	TYR	2.8
1	B	316	VAL	2.8
1	C	84	TRP	2.7
1	C	315	ALA	2.7
1	B	189	PRO	2.7
1	C	143	VAL	2.6
1	B	569	THR	2.5
1	C	388	GLU	2.5
1	C	453	GLN	2.4
1	C	146	TRP	2.4
1	A	85	GLY	2.4
1	C	316	VAL	2.4
1	B	87	LEU	2.3
1	C	101	LEU	2.3
1	C	602	ASP	2.3
1	A	611	MET	2.2
1	C	85	GLY	2.2

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Mol	Chain	Res	Type	RSRZ
1	B	88	THR	2.2
1	C	87	LEU	2.2
1	C	129	ALA	2.1
1	A	286	HIS	2.1
1	C	530	ILE	2.1
1	C	138	PHE	2.1
1	C	284	VAL	2.1
1	C	252	LEU	2.0
1	C	283	ASP	2.0
1	B	90	PRO	2.0
1	C	151	LEU	2.0

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

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

6.3 Carbohydrates [i](#)

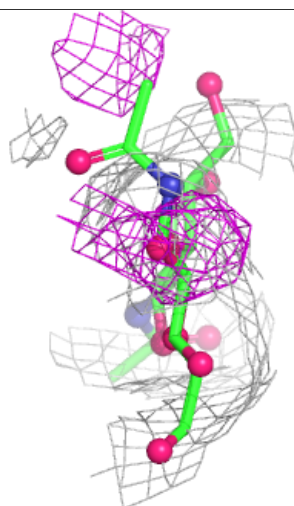
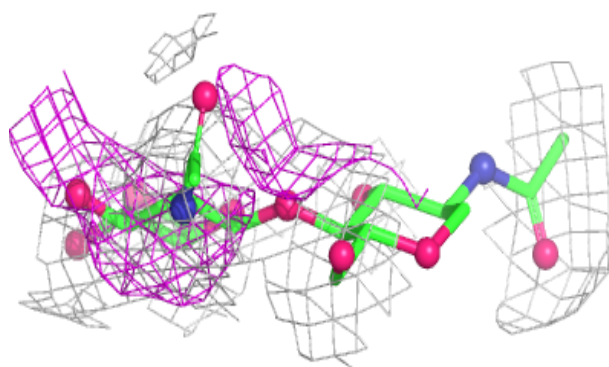
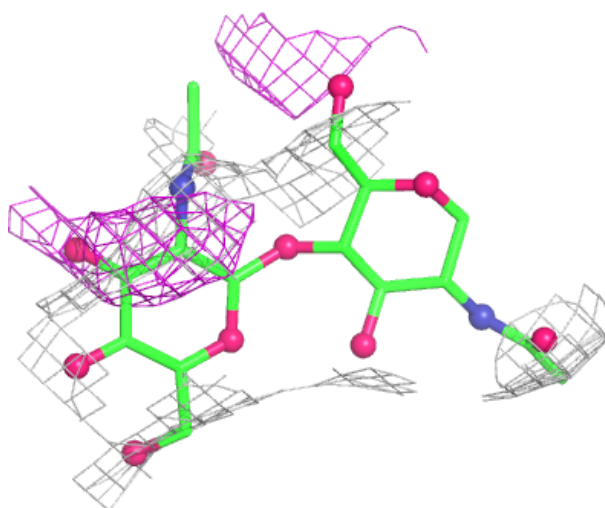
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q<0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
3	BMA	H	3	11/12	0.25	0.12	265,265,265,265	0
2	NAG	F	1	14/15	0.49	0.15	250,267,280,295	0
2	NAG	G	1	14/15	0.60	0.12	248,266,271,274	0
2	NAG	E	1	14/15	0.63	0.11	219,236,241,244	0
2	NAG	D	1	14/15	0.64	0.12	213,229,242,258	0
2	NAG	F	2	14/15	0.70	0.15	252,267,282,291	0
2	NAG	D	2	14/15	0.77	0.11	220,236,251,260	0
3	NAG	H	2	14/15	0.82	0.10	222,240,249,250	0
2	NAG	G	2	14/15	0.84	0.10	266,286,292,293	0
2	NAG	I	2	14/15	0.87	0.10	229,247,256,257	0
2	NAG	E	2	14/15	0.92	0.09	250,270,275,276	0
2	NAG	I	1	14/15	0.94	0.11	206,218,232,235	0
3	NAG	H	1	14/15	0.97	0.07	177,189,203,205	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.

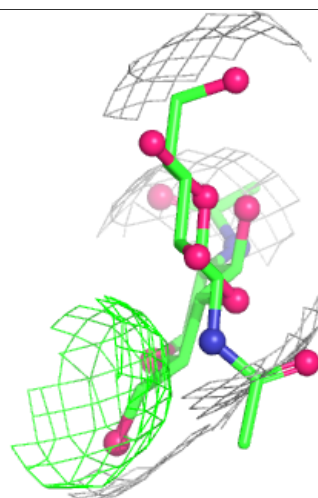
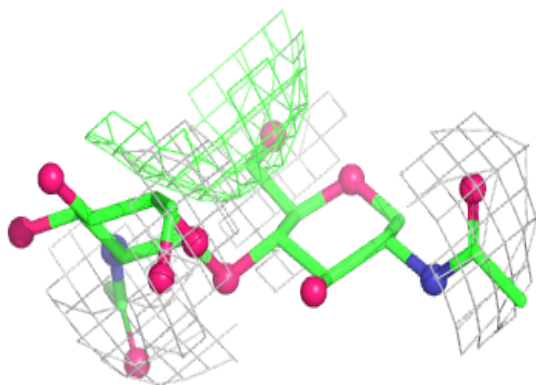
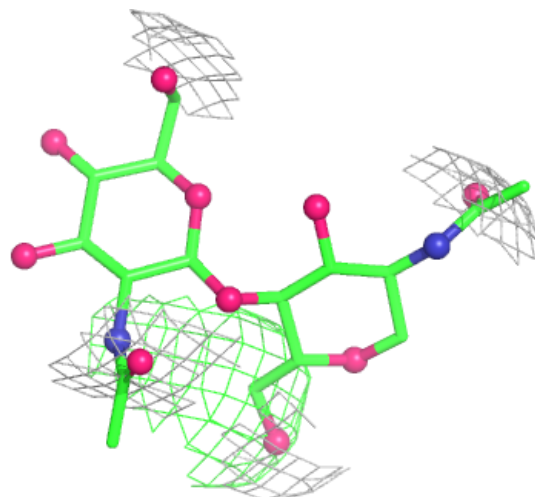
Electron density around Chain D:

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



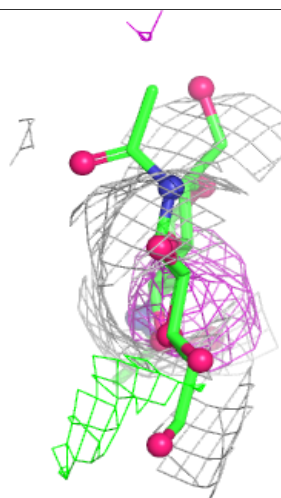
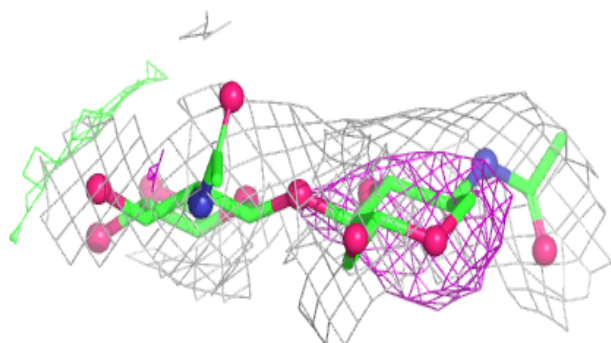
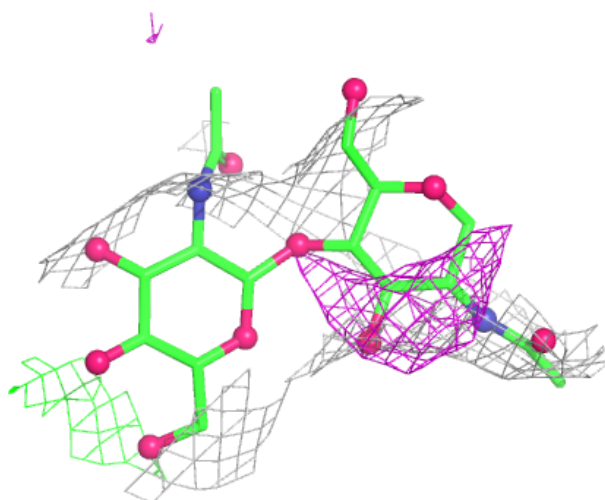
Electron density around Chain E:

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



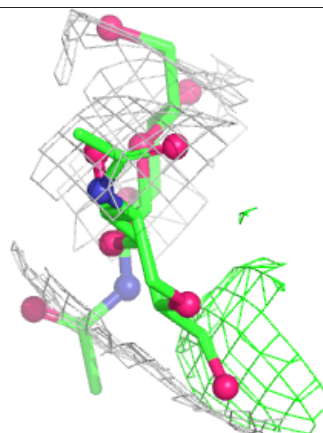
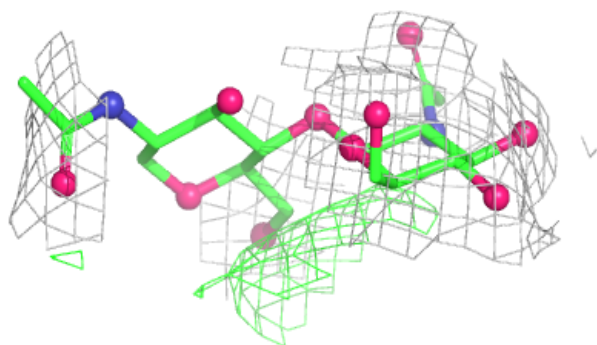
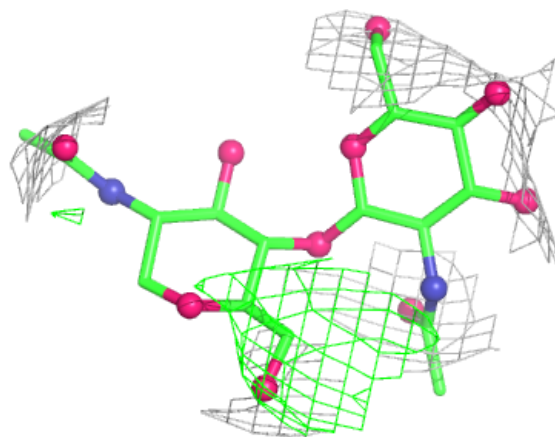
Electron density around Chain F:

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

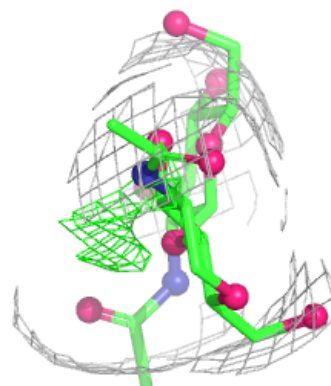
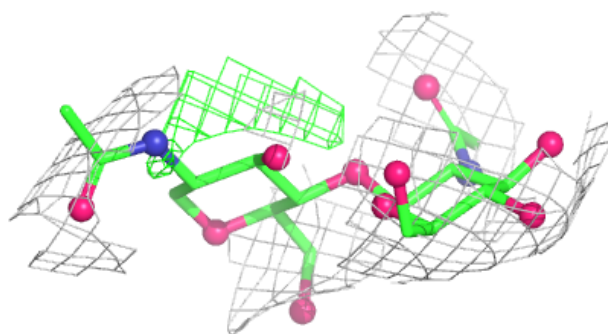
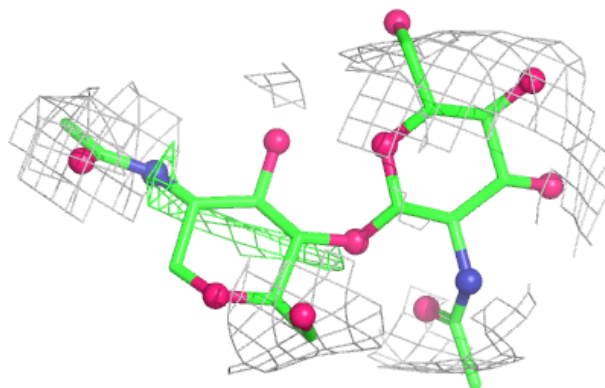


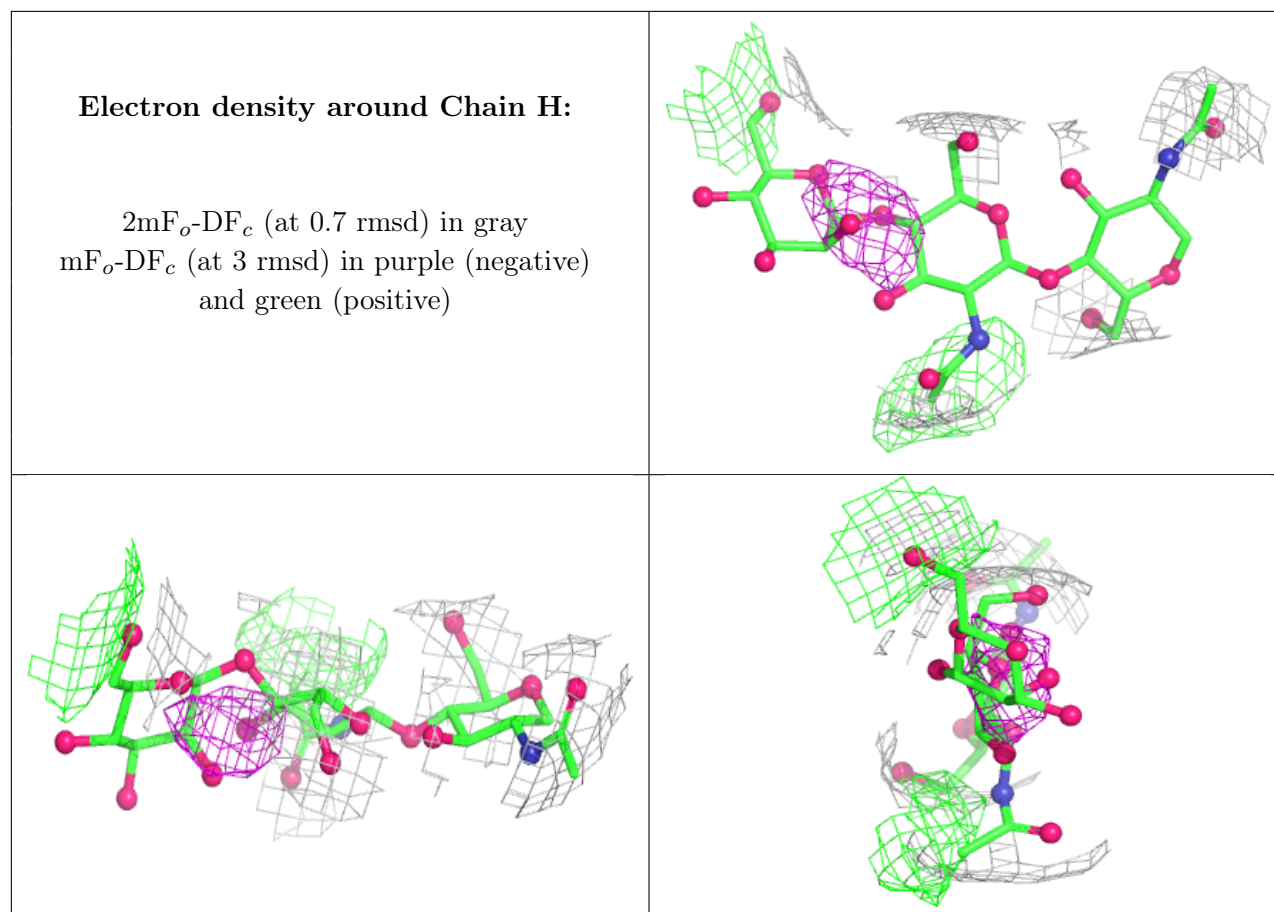
Electron density around Chain G:

$2mF_o-DF_c$ (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
5	BMA	A	707	11/12	0.23	0.10	292,292,292,292	0
4	NAG	C	703	14/15	0.28	0.16	314,332,337,340	0
4	NAG	C	702	14/15	0.40	0.13	303,329,337,337	0
5	BMA	B	709	11/12	0.49	0.07	309,309,309,309	0
6	MAN	A	711	11/12	0.58	0.08	279,279,279,279	0
4	NAG	B	701	14/15	0.60	0.15	294,295,297,298	0
4	NAG	A	701	14/15	0.60	0.15	286,288,290,291	0
4	NAG	C	701	14/15	0.61	0.16	337,339,341,341	0
4	NAG	A	702	14/15	0.68	0.12	250,276,284,284	0
4	NAG	B	708	14/15	0.68	0.08	301,301,301,301	0
4	NAG	C	704	14/15	0.73	0.10	313,323,334,336	0
4	NAG	B	702	14/15	0.76	0.12	266,292,300,301	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
4	NAG	A	710	14/15	0.87	0.10	234,252,261,262	0
4	NAG	A	708	14/15	0.88	0.09	215,226,236,239	0
4	NAG	B	707	14/15	0.88	0.10	267,277,288,290	0
7	ZN	C	708	1/1	0.91	0.07	290,290,290,290	0
7	ZN	A	713	1/1	0.92	0.11	176,176,176,176	0
4	NAG	A	709	14/15	0.93	0.08	179,192,206,208	0
7	ZN	C	707	1/1	0.94	0.07	297,297,297,297	0
7	ZN	A	712	1/1	0.95	0.10	175,175,175,175	0
7	ZN	B	714	1/1	0.95	0.07	189,189,189,189	0
7	ZN	B	713	1/1	0.97	0.07	202,202,202,202	0

6.5 Other polymers [i](#)

There are no such residues in this entry.