



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 7, 2026 – 01:31 AM UTC

PDB ID : 7CZF / pdb_00007czf
Title : Crystal structure of Kaposi Sarcoma associated herpesvirus (KSHV) gHgL in complex with the ligand binding domian (LBD) of EphA2
Authors : Su, C.; Wu, L.L.; Song, H.; Chai, Y.; Qi, J.X.; Yan, J.H.; Gao, G.F.
Deposited on : 2020-09-08
Resolution : 3.20 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

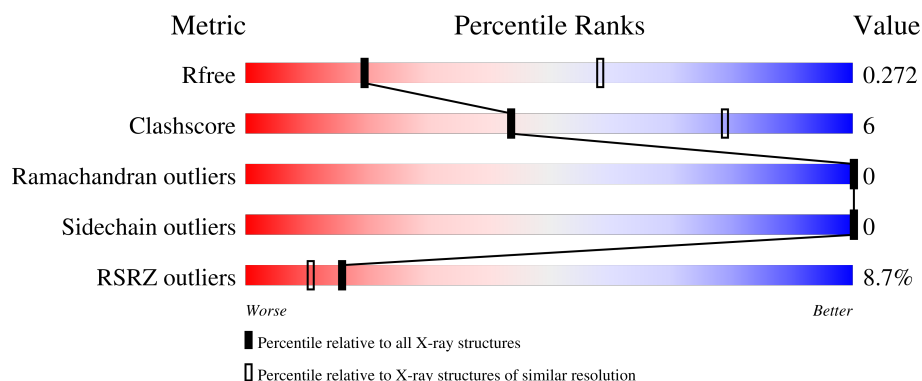
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.20 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.



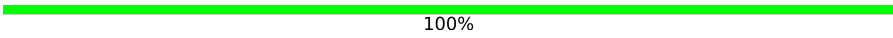
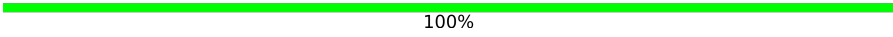
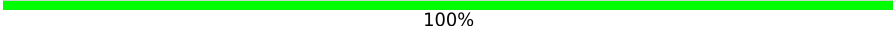
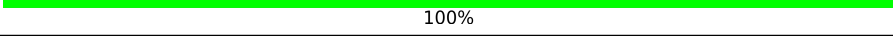
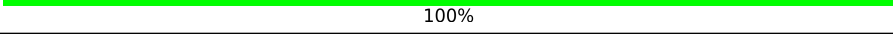
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	1466 (3.20-3.20)
Clashscore	190562	1573 (3.20-3.20)
Ramachandran outliers	187476	1548 (3.20-3.20)
Sidechain outliers	187428	1547 (3.20-3.20)
RSRZ outliers	180081	1466 (3.20-3.20)

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	185	11% 76% 16% 8%
1	D	185	21% 73% 12% 15%
2	B	677	5% 79% 16% 5%
2	E	677	6% 78% 17% 5%
3	C	157	6% 60% 13% 27%

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Mol	Chain	Length	Quality of chain
3	F	157	
4	K	2	
4	L	2	
4	N	2	
4	a	2	
5	S	6	
6	H	3	

2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 14901 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 Ephrin type-A receptor 2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	170	Total	C	N	O	S	0	0	0
			1376	883	231	252	10			
1	D	158	Total	C	N	O	S	0	0	0
			1271	817	208	237	9			

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	207	HIS	-	expression tag	UNP P29317
A	208	HIS	-	expression tag	UNP P29317
A	209	HIS	-	expression tag	UNP P29317
A	210	HIS	-	expression tag	UNP P29317
A	211	HIS	-	expression tag	UNP P29317
A	212	HIS	-	expression tag	UNP P29317
D	207	HIS	-	expression tag	UNP P29317
D	208	HIS	-	expression tag	UNP P29317
D	209	HIS	-	expression tag	UNP P29317
D	210	HIS	-	expression tag	UNP P29317
D	211	HIS	-	expression tag	UNP P29317
D	212	HIS	-	expression tag	UNP P29317

- Molecule 2 is a protein called Envelope glycoprotein H.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	644	Total	C	N	O	S	0	0	0
			5061	3237	851	945	28			
2	E	642	Total	C	N	O	S	0	0	0
			5049	3229	849	943	28			

- Molecule 3 is a protein called Envelope glycoprotein L.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	114	Total	C	N	O	S	0	0	0
			874	554	156	159	5			
3	F	115	Total	C	N	O	S	0	0	0
			882	562	157	158	5			

There are 20 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	17	ASP	-	expression tag	UNP Q76RG7
C	18	GLY	-	expression tag	UNP Q76RG7
C	19	ILE	-	expression tag	UNP Q76RG7
C	20	GLN	-	expression tag	UNP Q76RG7
C	168	HIS	-	expression tag	UNP Q76RG7
C	169	HIS	-	expression tag	UNP Q76RG7
C	170	HIS	-	expression tag	UNP Q76RG7
C	171	HIS	-	expression tag	UNP Q76RG7
C	172	HIS	-	expression tag	UNP Q76RG7
C	173	HIS	-	expression tag	UNP Q76RG7
F	17	ASP	-	expression tag	UNP Q76RG7
F	18	GLY	-	expression tag	UNP Q76RG7
F	19	ILE	-	expression tag	UNP Q76RG7
F	20	GLN	-	expression tag	UNP Q76RG7
F	168	HIS	-	expression tag	UNP Q76RG7
F	169	HIS	-	expression tag	UNP Q76RG7
F	170	HIS	-	expression tag	UNP Q76RG7
F	171	HIS	-	expression tag	UNP Q76RG7
F	172	HIS	-	expression tag	UNP Q76RG7
F	173	HIS	-	expression tag	UNP Q76RG7

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



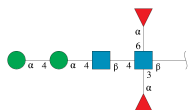
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
4	L	2	Total	C	N	O	0	0	0
			28	16	2	10			
4	a	2	Total	C	N	O	0	0	0
			28	16	2	10			
4	K	2	Total	C	N	O	0	0	0
			28	16	2	10			

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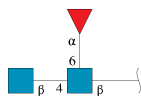
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
4	N	2	Total	C	N	O	0	0	0
			28	16	2	10			

- Molecule 5 is an oligosaccharide called alpha-D-mannopyranose-(1-4)-alpha-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-[alpha-L-fucopyranose-(1-3)][alpha-L-fucopyranose-(1-6)]2-acetamido-2-deoxy-beta-D-glucopyranose.



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
5	S	6	Total	C	N	O	0	0	0
			70	40	2	28			

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



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
6	H	3	Total	C	N	O	0	0	0
			38	22	2	14			

- Molecule 7 is 2-acetamido-2-deoxy-beta-D-glucopyranose (CCD ID: NAG) (formula: C₈H₁₅NO₆).

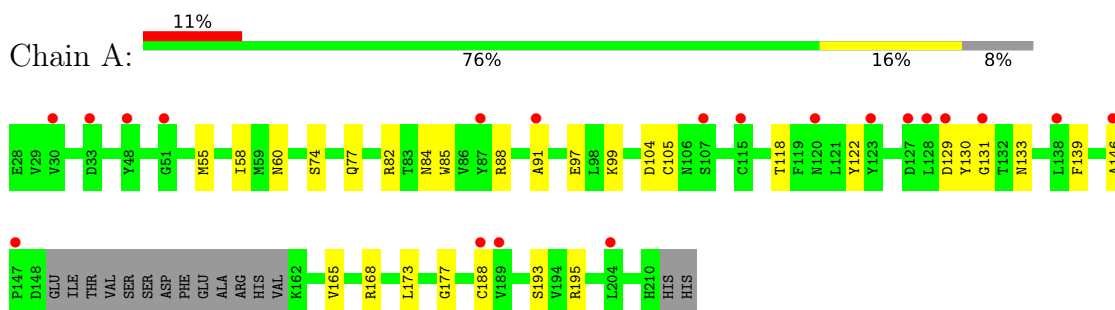


Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
7	B	1	Total	C	N	O	0	0
			14	8	1	5		
7	B	1	Total	C	N	O	0	0
			14	8	1	5		
7	B	1	Total	C	N	O	0	0
			14	8	1	5		
7	B	1	Total	C	N	O	0	0
			14	8	1	5		
7	B	1	Total	C	N	O	0	0
			14	8	1	5		
7	C	1	Total	C	N	O	0	0
			14	8	1	5		
7	E	1	Total	C	N	O	0	0
			14	8	1	5		
7	E	1	Total	C	N	O	0	0
			14	8	1	5		
7	E	1	Total	C	N	O	0	0
			14	8	1	5		
7	E	1	Total	C	N	O	0	0
			14	8	1	5		
7	F	1	Total	C	N	O	0	0
			14	8	1	5		

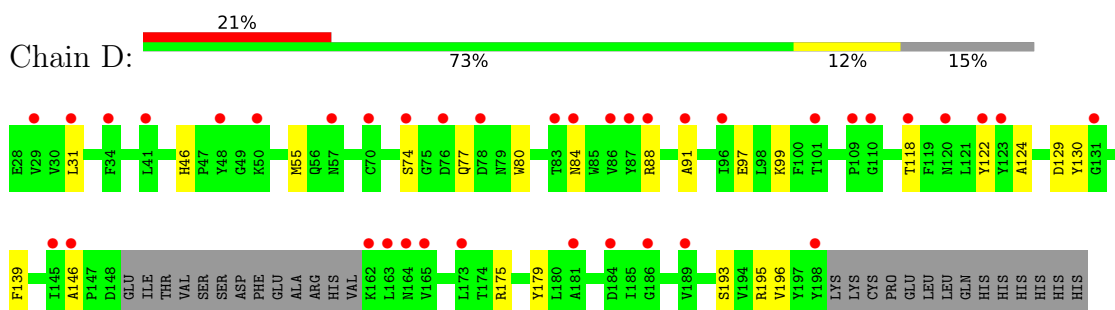
3 Residue-property plots [i](#)

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.

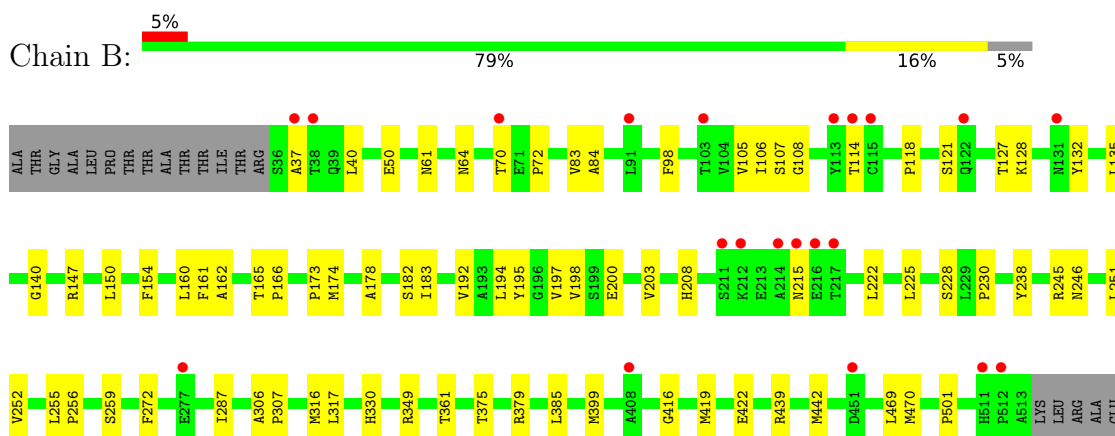
• Molecule 1: Ephrin type-A receptor 2



• Molecule 1: Ephrin type-A receptor 2



• Molecule 2: Envelope glycoprotein H



Chain L:  100%

MAG1
MAG2

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

Chain a:  100%

MAG1
MAG2

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

Chain K:  100%

MAG1
MAG2

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

Chain N:  100%

MAG1
MAG2

- Molecule 5: alpha-D-mannopyranose-(1-4)-alpha-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-[alpha-L-fucopyranose-(1-3)][alpha-L-fucopyranose-(1-6)]2-acetamido-2-deoxy-beta-D-glucopyranose

Chain S:  67% 33%

MAG1
MAG2
MAN3
MAN4
FUC5
FUC6

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

Chain H:  100%

MAG1
MAG2
FUC3

4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	80.80Å 153.20Å 275.23Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	31.65 – 3.20 31.65 – 3.20	Depositor EDS
% Data completeness (in resolution range)	78.7 (31.65-3.20) 78.7 (31.65-3.20)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.96 (at 3.18Å)	Xtriage
Refinement program	PHENIX 1.16_3549	Depositor
R, R_{free}	0.239 , 0.272 0.240 , 0.272	Depositor DCC
R_{free} test set	2201 reflections (3.83%)	wwPDB-VP
Wilson B-factor (Å ²)	54.1	Xtriage
Anisotropy	0.303	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.27 , 35.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.87	EDS
Total number of atoms	14901	wwPDB-VP
Average B, all atoms (Å ²)	64.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.15% 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: FUC, MAN, NAG

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	0.07	0/1412	0.25	0/1911
1	D	0.08	0/1302	0.24	0/1763
2	B	0.09	0/5172	0.31	0/7033
2	E	0.09	0/5159	0.30	0/7014
3	C	0.12	0/893	0.31	0/1218
3	F	0.09	0/901	0.30	0/1229
All	All	0.09	0/14839	0.30	0/20168

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

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	1376	0	1319	20	0
1	D	1271	0	1217	14	0
2	B	5061	0	5031	62	0
2	E	5049	0	5019	69	0
3	C	874	0	858	18	0
3	F	882	0	876	11	0
4	K	28	0	25	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	L	28	0	25	0	0
4	N	28	0	25	0	0
4	a	28	0	25	0	0
5	S	70	0	61	0	0
6	H	38	0	34	0	0
7	B	70	0	65	1	0
7	C	14	0	13	0	0
7	E	70	0	65	1	0
7	F	14	0	13	0	0
All	All	14901	0	14671	179	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 6.

All (179) 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)
2:E:674:ASP:O	2:E:677:ASN:ND2	2.22	0.73
2:B:627:ILE:HG13	2:B:696:GLY:HA3	1.70	0.72
2:B:105:VAL:HG22	2:B:245:ARG:HB3	1.73	0.70
2:B:606:LYS:HD2	2:B:610:SER:HB2	1.73	0.69
2:E:470:MET:HE3	2:E:657:THR:HG21	1.75	0.69
2:B:553:PRO:HB2	2:B:607:PRO:HB3	1.75	0.68
3:C:57:VAL:CG2	3:C:59:ILE:HD12	2.23	0.67
2:B:552:ILE:HB	2:B:555:ILE:HB	1.76	0.66
2:E:636:LEU:HD12	2:E:662:GLN:HG3	1.78	0.66
1:A:84:ASN:HD22	1:D:129:ASP:HB3	1.61	0.65
2:E:578:ILE:HG12	2:E:642:MET:HE1	1.79	0.65
2:B:222:LEU:HB2	2:B:252:VAL:HB	1.80	0.63
2:B:127:THR:HG22	2:B:128:LYS:H	1.62	0.63
1:A:122:TYR:HB3	1:A:139:PHE:HB3	1.81	0.62
2:B:70:THR:HG22	2:B:72:PRO:HD2	1.82	0.62
2:B:630:PRO:O	2:B:662:GLN:NE2	2.23	0.62
2:E:129:GLY:O	2:E:131:ASN:ND2	2.31	0.61
2:B:259:SER:HA	7:B:702:NAG:H83	1.81	0.61
3:C:57:VAL:HG23	3:C:59:ILE:HD12	1.82	0.61
2:B:147:ARG:HH21	2:B:150:LEU:HD13	1.65	0.61
1:D:74:SER:OG	1:D:77:GLN:OE1	2.19	0.60
1:D:46:HIS:HB3	1:D:80:TRP:HB2	1.84	0.60
2:E:241:LEU:HD13	2:E:254:ILE:HG12	1.84	0.59
1:D:122:TYR:HB3	1:D:139:PHE:HB3	1.83	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:E:44:ARG:HD3	2:E:44:ARG:H	1.66	0.59
2:E:222:LEU:HD12	2:E:229:LEU:HD22	1.85	0.59
2:B:203:VAL:HG21	2:B:317:LEU:HB3	1.85	0.57
2:E:186:GLY:HA3	2:E:191:GLY:HA3	1.87	0.57
2:E:469:LEU:HD13	2:E:570:PRO:HD2	1.85	0.57
2:E:552:ILE:HB	2:E:555:ILE:HB	1.85	0.57
2:E:660:ARG:O	2:E:664:ASN:ND2	2.34	0.57
1:A:97:GLU:HB3	1:A:195:ARG:HB2	1.87	0.56
1:A:130:TYR:HB3	1:A:133:ASN:HB2	1.87	0.56
2:E:49:ILE:HD13	3:F:46:ILE:HG13	1.87	0.56
1:A:74:SER:OG	1:A:77:GLN:OE1	2.22	0.56
2:B:686:ARG:HG3	2:B:688:ASN:HB2	1.88	0.56
2:E:65:LEU:HB3	2:E:74:LEU:HD11	1.88	0.55
1:D:124:ALA:HB2	1:D:139:PHE:HA	1.87	0.55
3:C:81:ARG:HG2	3:C:106:LEU:HD22	1.89	0.55
1:A:88:ARG:HG3	1:A:91:ALA:HB3	1.88	0.55
1:D:97:GLU:OE1	1:D:195:ARG:NH1	2.39	0.55
1:D:88:ARG:HG3	1:D:91:ALA:HB3	1.89	0.55
3:C:57:VAL:HG21	3:C:59:ILE:HD12	1.88	0.54
2:B:50:GLU:HG3	3:C:19:ILE:HG13	1.90	0.54
2:B:583:LEU:HD23	2:B:605:ILE:HD11	1.91	0.53
1:A:55:MET:HG2	3:C:70:THR:HB	1.90	0.53
2:E:571:LEU:HD21	2:E:640:VAL:HG23	1.90	0.53
3:C:39:LEU:HD23	3:C:39:LEU:H	1.74	0.53
2:B:173:PRO:O	3:C:90:ARG:NH2	2.42	0.52
2:B:106:ILE:HG13	2:B:107:SER:H	1.75	0.52
2:E:622:PRO:HB2	2:E:692:VAL:HG23	1.90	0.52
2:B:140:GLY:HA3	2:B:399:MET:HE1	1.92	0.52
1:A:85:TRP:HE1	1:A:177:GLY:HA3	1.75	0.51
2:B:422:GLU:HG3	2:B:470:MET:HE1	1.92	0.51
2:B:660:ARG:O	2:B:664:ASN:ND2	2.30	0.51
2:E:554:ALA:HB2	2:E:607:PRO:HA	1.92	0.51
2:E:606:LYS:HD2	2:E:610:SER:HB2	1.93	0.51
2:E:684:TRP:HB3	2:E:692:VAL:HG12	1.93	0.51
2:E:686:ARG:HH21	2:E:688:ASN:HD22	1.58	0.51
2:B:230:PRO:O	2:B:238:TYR:OH	2.22	0.51
2:E:132:TYR:HB3	2:E:375:THR:HG21	1.92	0.51
2:E:222:LEU:HB2	2:E:252:VAL:HB	1.93	0.50
1:A:129:ASP:HB3	1:D:84:ASN:HD22	1.77	0.50
2:B:571:LEU:HD21	2:B:640:VAL:HG23	1.92	0.50
2:B:564:LYS:HB3	2:B:580:SER:HB2	1.94	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:F:37:LEU:HD12	3:F:38:PRO:HD2	1.92	0.50
2:E:383:THR:HA	2:E:419:MET:HE1	1.94	0.50
2:E:641:ILE:HD12	2:E:665:LEU:HD23	1.92	0.50
2:E:122:GLN:HB3	2:E:124:TYR:HE1	1.77	0.50
2:B:174:MET:HE1	2:B:230:PRO:HA	1.93	0.50
3:C:46:ILE:HG22	3:C:46:ILE:O	2.12	0.50
2:E:84:ALA:HB2	3:F:37:LEU:HD13	1.94	0.50
2:E:327:ILE:HG23	2:E:333:VAL:HG11	1.94	0.49
2:E:167:GLU:HG3	2:E:220:HIS:NE2	2.28	0.49
1:A:97:GLU:OE1	1:A:195:ARG:NH1	2.40	0.49
2:E:448:VAL:O	2:E:452:SER:OG	2.22	0.49
2:B:225:LEU:HB2	2:B:228:SER:HB2	1.95	0.49
2:E:146:ALA:HB3	2:E:331:ARG:HA	1.94	0.48
2:B:287:ILE:HG21	2:B:316:MET:HG3	1.95	0.48
2:B:349:ARG:HD2	2:B:361:THR:HG21	1.95	0.48
1:A:58:ILE:HG22	3:C:30:GLN:HG3	1.94	0.48
2:E:388:LEU:HB3	2:E:392:SER:HB2	1.95	0.48
2:E:390:ARG:NH1	2:E:438:GLU:OE2	2.47	0.48
2:E:571:LEU:HB2	2:E:574:VAL:HG13	1.95	0.48
2:E:286:THR:O	2:E:290:ARG:HG2	2.14	0.48
2:E:354:LEU:HD21	2:E:544:LEU:HD12	1.96	0.48
2:E:106:ILE:HG22	2:E:107:SER:H	1.79	0.48
2:E:634:CYS:HA	2:E:662:GLN:HE22	1.79	0.48
2:B:132:TYR:HB3	2:B:375:THR:HG21	1.96	0.47
2:B:215:ASN:O	2:B:215:ASN:ND2	2.47	0.47
2:B:667:LEU:O	2:B:670:SER:OG	2.27	0.47
2:E:150:LEU:HD11	2:E:328:ALA:HB2	1.95	0.47
2:E:230:PRO:O	2:E:238:TYR:OH	2.31	0.47
1:A:60:ASN:N	3:C:30:GLN:OE1	2.48	0.47
2:B:566:ILE:HG21	2:B:650:LEU:HD13	1.95	0.47
2:E:106:ILE:HG21	2:E:273:LEU:HD21	1.97	0.47
2:B:192:VAL:HG11	2:B:306:ALA:HB1	1.96	0.47
2:E:183:ILE:N	2:E:194:LEU:O	2.32	0.47
2:E:694:ILE:HG22	2:E:698:TYR:HE1	1.80	0.47
2:E:259:SER:HA	7:E:702:NAG:H83	1.95	0.47
1:D:88:ARG:NE	1:D:175:ARG:O	2.48	0.47
1:D:118:THR:HG22	1:D:146:ALA:HA	1.96	0.47
2:E:167:GLU:OE2	2:E:206:THR:HG21	2.15	0.46
2:E:117:VAL:HG21	2:E:123:THR:HG22	1.97	0.46
1:A:188:CYS:SG	3:C:23:ALA:HA	2.55	0.46
1:A:118:THR:HG22	1:A:146:ALA:HA	1.97	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:469:LEU:HD13	2:B:570:PRO:HD2	1.97	0.46
2:B:108:GLY:HA2	2:B:246:ASN:HB2	1.96	0.46
2:B:160:LEU:HD11	2:B:197:VAL:HG11	1.96	0.46
3:F:91:LEU:O	3:F:95:ALA:N	2.48	0.46
2:B:251:LEU:HD22	2:B:272:PHE:HB2	1.98	0.45
2:E:185:ILE:HD12	2:E:303:GLN:HA	1.98	0.45
2:B:469:LEU:HD11	2:B:569:VAL:HG13	1.99	0.45
2:B:641:ILE:HD12	2:B:665:LEU:HD23	1.98	0.45
2:E:78:TRP:CD2	3:F:40:PHE:HB2	2.52	0.45
2:E:246:ASN:HB2	2:E:272:PHE:CZ	2.52	0.45
2:B:84:ALA:HB2	3:C:37:LEU:HD13	1.99	0.45
3:C:75:VAL:HG11	3:C:80:LEU:HD22	1.98	0.45
2:E:205:VAL:HG22	2:E:317:LEU:HD22	1.98	0.45
2:E:354:LEU:HB3	2:E:540:LEU:HD13	2.00	0.44
2:E:572:PRO:O	2:E:573:HIS:ND1	2.50	0.44
2:E:109:ASP:N	2:E:246:ASN:OD1	2.50	0.44
2:B:162:ALA:O	2:B:165:THR:OG1	2.28	0.44
3:C:22:VAL:HG13	3:C:63:ARG:HH21	1.81	0.44
3:F:118:ASN:OD1	3:F:118:ASN:N	2.45	0.44
2:B:37:ALA:HB1	2:B:40:LEU:HB2	1.99	0.44
2:B:98:PHE:HB3	2:B:255:LEU:HD12	1.99	0.44
1:A:104:ASP:OD1	1:A:105:CYS:N	2.50	0.44
2:B:439:ARG:HA	2:B:442:MET:HE2	1.98	0.44
1:D:55:MET:HG2	3:F:70:THR:HB	2.00	0.44
2:B:183:ILE:N	2:B:194:LEU:O	2.33	0.44
2:E:564:LYS:HB3	2:E:580:SER:HB2	2.00	0.44
3:C:101:ARG:HA	3:C:104:TRP:HB2	1.98	0.44
2:E:335:VAL:HG13	2:E:369:LEU:HD11	2.00	0.44
2:E:493:LEU:HD23	2:E:493:LEU:HA	1.78	0.43
2:B:632:ARG:H	2:B:632:ARG:HD3	1.84	0.43
1:A:88:ARG:HD3	1:A:173:LEU:HD13	1.99	0.43
2:B:573:HIS:HA	2:B:607:PRO:HG2	2.01	0.43
2:B:83:VAL:HB	3:C:75:VAL:HG13	2.00	0.42
2:E:428:TYR:CD2	2:E:474:MET:HG2	2.54	0.42
3:C:61:LYS:HE2	3:C:72:GLU:HB3	2.01	0.42
2:E:653:LEU:O	2:E:654:MET:HG2	2.20	0.42
2:E:78:TRP:CG	3:F:40:PHE:HB2	2.54	0.42
2:E:174:MET:HE1	2:E:230:PRO:HA	2.01	0.42
3:F:104:TRP:O	3:F:108:ARG:HG2	2.18	0.42
1:A:99:LYS:HA	1:A:165:VAL:HG22	2.02	0.42
2:B:114:THR:HB	2:B:330:HIS:NE2	2.34	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:118:PRO:O	2:B:121:SER:HB3	2.19	0.42
2:B:154:PHE:CE2	2:B:200:GLU:HB3	2.54	0.42
2:E:313:LEU:O	2:E:317:LEU:HD13	2.19	0.42
1:D:99:LYS:HB2	1:D:193:SER:HB3	2.00	0.42
2:B:256:PRO:HD2	2:B:259:SER:HB2	2.02	0.42
2:B:419:MET:HE3	2:B:419:MET:HB3	1.88	0.42
2:E:563:ASP:OD1	2:E:563:ASP:N	2.51	0.42
3:F:57:VAL:HG12	3:F:121:THR:HA	2.02	0.42
1:D:130:TYR:N	1:D:179:TYR:OH	2.49	0.42
2:B:182:SER:HA	2:B:195:TYR:HA	2.01	0.42
1:A:82:ARG:NH1	1:A:131:GLY:O	2.46	0.42
2:B:61:ASN:HB3	2:B:64:ASN:HB3	2.01	0.42
1:D:31:LEU:N	1:D:196:VAL:O	2.44	0.41
2:E:212:LYS:HA	2:E:212:LYS:HD2	1.92	0.41
2:E:401:GLN:HA	2:E:404:ALA:HB3	2.02	0.41
2:E:152:ASP:HB3	2:E:366:HIS:HA	2.01	0.41
2:E:473:SER:HA	2:E:653:LEU:O	2.21	0.41
3:F:43:VAL:HG11	3:F:62:LEU:HD22	2.02	0.41
2:B:166:PRO:HG3	2:B:208:HIS:CG	2.56	0.41
2:B:178:ALA:HB1	2:B:198:VAL:O	2.21	0.41
2:B:135:LEU:O	2:B:379:ARG:NH2	2.53	0.41
2:B:306:ALA:N	2:B:307:PRO:HD2	2.35	0.41
2:E:37:ALA:HB1	2:E:40:LEU:HB2	2.03	0.41
2:B:161:PHE:HB2	2:B:173:PRO:HA	2.03	0.41
2:B:571:LEU:HB2	2:B:574:VAL:HG13	2.03	0.41
1:A:168:ARG:HD3	1:A:168:ARG:HA	1.96	0.41
2:E:243:PHE:CE1	2:E:253:ALA:HB3	2.56	0.41
2:E:306:ALA:N	2:E:307:PRO:HD2	2.36	0.41
1:A:99:LYS:HB2	1:A:193:SER:HB3	2.03	0.40
2:B:385:LEU:HD23	2:B:416:GLY:HA3	2.04	0.40
2:E:96:SER:O	2:E:257:LYS:N	2.42	0.40
2:B:442:MET:HE1	2:B:501:PRO:HD3	2.04	0.40

There are no symmetry-related clashes.

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

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

of similar resolution.

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	166/185 (90%)	166 (100%)	0	0	100	100
1	D	154/185 (83%)	154 (100%)	0	0	100	100
2	B	640/677 (94%)	640 (100%)	0	0	100	100
2	E	638/677 (94%)	636 (100%)	2 (0%)	0	100	100
3	C	112/157 (71%)	112 (100%)	0	0	100	100
3	F	113/157 (72%)	113 (100%)	0	0	100	100
All	All	1823/2038 (90%)	1821 (100%)	2 (0%)	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	147/161 (91%)	147 (100%)	0	100	100
1	D	135/161 (84%)	135 (100%)	0	100	100
2	B	562/586 (96%)	562 (100%)	0	100	100
2	E	561/586 (96%)	561 (100%)	0	100	100
3	C	95/129 (74%)	95 (100%)	0	100	100
3	F	96/129 (74%)	96 (100%)	0	100	100
All	All	1596/1752 (91%)	1596 (100%)	0	100	100

There are no protein residues with a non-rotameric sidechain to report.

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (14) such sidechains are listed below:

Mol	Chain	Res	Type
2	B	64	ASN

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Mol	Chain	Res	Type
2	B	215	ASN
2	B	288	GLN
2	B	303	GLN
2	B	679	HIS
3	C	107	GLN
3	C	130	ASN
1	D	120	ASN
2	E	42	ASN
2	E	477	ASN
2	E	494	ASN
2	E	620	HIS
3	F	44	HIS
3	F	130	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

17 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	$\# Z > 2$
6	NAG	H	1	6,2	14,14,15	0.29	0	17,19,21	0.46	0
6	NAG	H	2	6	14,14,15	0.22	0	17,19,21	0.41	0
6	FUC	H	3	6	10,10,11	0.77	0	14,14,16	0.89	0
4	NAG	K	1	4,2	14,14,15	0.25	0	17,19,21	0.48	0
4	NAG	K	2	4	14,14,15	0.21	0	17,19,21	0.42	0
4	NAG	L	1	4,2	14,14,15	0.28	0	17,19,21	0.51	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
4	NAG	L	2	4	14,14,15	0.34	0	17,19,21	0.49	0
4	NAG	N	1	4,2	14,14,15	0.25	0	17,19,21	0.47	0
4	NAG	N	2	4	14,14,15	0.31	0	17,19,21	0.39	0
5	NAG	S	1	2,5	14,14,15	0.21	0	17,19,21	0.42	0
5	NAG	S	2	5	14,14,15	0.22	0	17,19,21	0.45	0
5	MAN	S	3	5	11,11,12	0.68	0	15,15,17	0.96	2 (13%)
5	MAN	S	4	5	11,11,12	0.72	0	15,15,17	0.99	2 (13%)
5	FUC	S	5	5	10,10,11	0.75	0	14,14,16	0.72	0
5	FUC	S	6	5	10,10,11	0.74	0	14,14,16	0.75	0
4	NAG	a	1	4,2	14,14,15	0.25	0	17,19,21	0.47	0
4	NAG	a	2	4	14,14,15	0.23	0	17,19,21	0.42	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
6	NAG	H	1	6,2	-	2/6/23/26	0/1/1/1
6	NAG	H	2	6	-	0/6/23/26	0/1/1/1
6	FUC	H	3	6	-	-	0/1/1/1
4	NAG	K	1	4,2	-	2/6/23/26	0/1/1/1
4	NAG	K	2	4	-	2/6/23/26	0/1/1/1
4	NAG	L	1	4,2	-	0/6/23/26	0/1/1/1
4	NAG	L	2	4	-	3/6/23/26	0/1/1/1
4	NAG	N	1	4,2	-	1/6/23/26	0/1/1/1
4	NAG	N	2	4	-	2/6/23/26	0/1/1/1
5	NAG	S	1	2,5	-	2/6/23/26	0/1/1/1
5	NAG	S	2	5	-	0/6/23/26	0/1/1/1
5	MAN	S	3	5	-	2/2/19/22	0/1/1/1
5	MAN	S	4	5	-	2/2/19/22	0/1/1/1
5	FUC	S	5	5	-	-	0/1/1/1
5	FUC	S	6	5	-	-	0/1/1/1
4	NAG	a	1	4,2	-	2/6/23/26	0/1/1/1
4	NAG	a	2	4	-	2/6/23/26	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(°)
5	S	3	MAN	C1-O5-C5	2.44	115.45	112.19
5	S	4	MAN	C1-O5-C5	2.40	115.40	112.19
5	S	3	MAN	O2-C2-C3	-2.16	105.67	110.15
5	S	4	MAN	O2-C2-C3	-2.14	105.71	110.15

There are no chirality outliers.

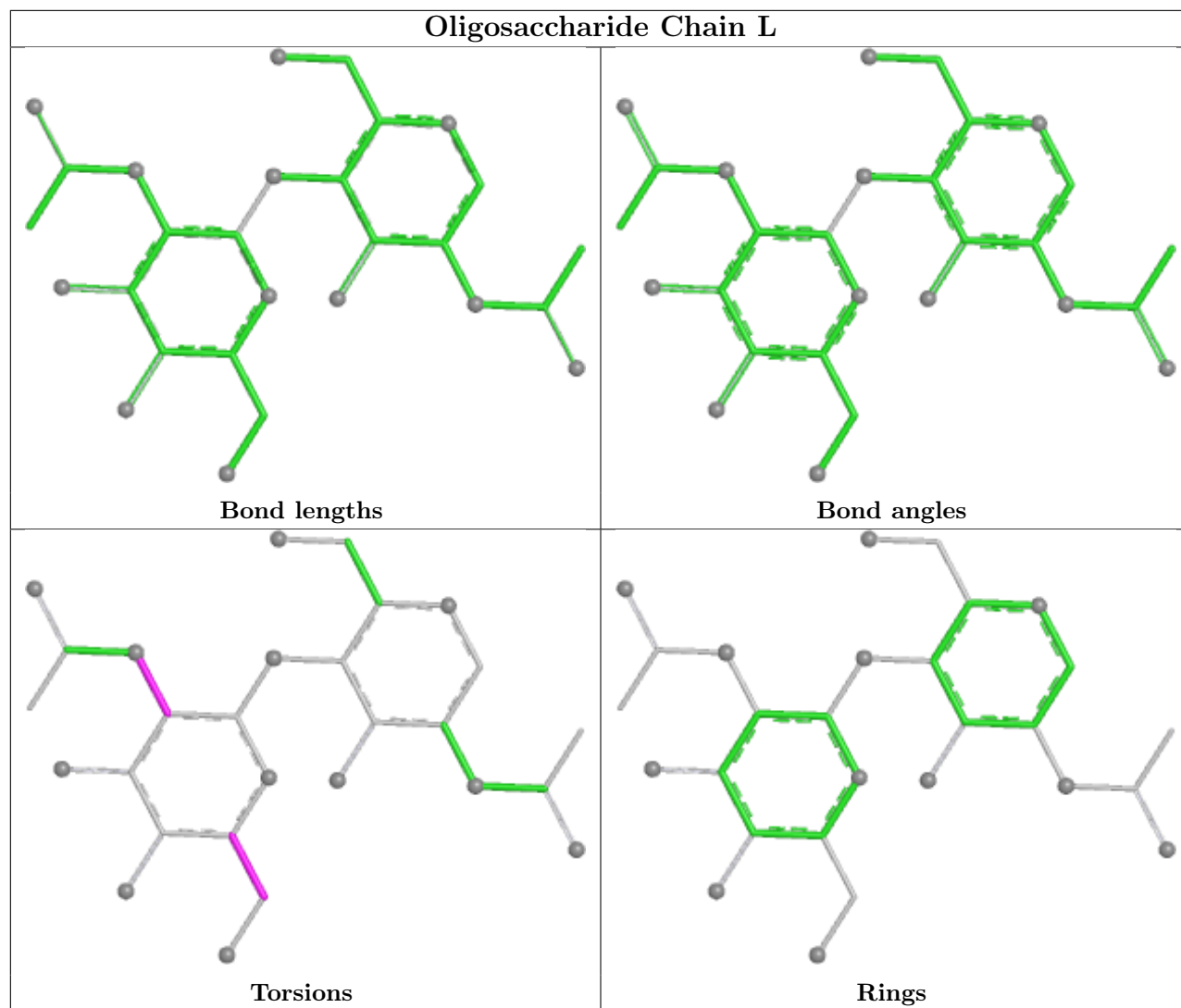
All (22) torsion outliers are listed below:

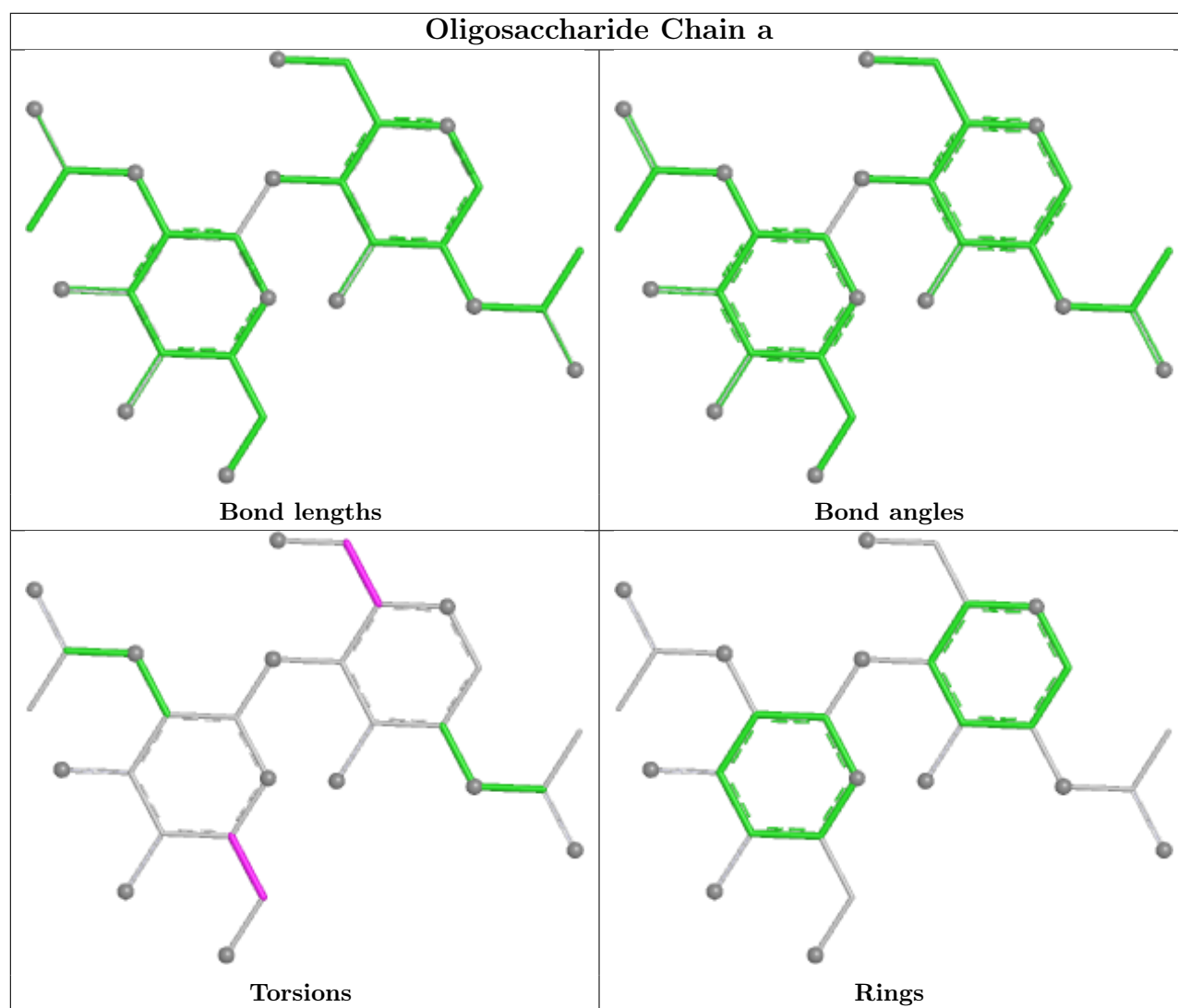
Mol	Chain	Res	Type	Atoms
4	L	2	NAG	O5-C5-C6-O6
5	S	1	NAG	O5-C5-C6-O6
4	K	1	NAG	O5-C5-C6-O6
4	L	2	NAG	C4-C5-C6-O6
5	S	1	NAG	C4-C5-C6-O6
4	K	2	NAG	C4-C5-C6-O6
6	H	1	NAG	O5-C5-C6-O6
4	a	2	NAG	C4-C5-C6-O6
4	K	1	NAG	C4-C5-C6-O6
4	K	2	NAG	O5-C5-C6-O6
4	N	2	NAG	C4-C5-C6-O6
4	a	2	NAG	O5-C5-C6-O6
4	a	1	NAG	C4-C5-C6-O6
4	N	2	NAG	O5-C5-C6-O6
6	H	1	NAG	C4-C5-C6-O6
4	a	1	NAG	O5-C5-C6-O6
5	S	3	MAN	O5-C5-C6-O6
5	S	4	MAN	C4-C5-C6-O6
4	N	1	NAG	C4-C5-C6-O6
5	S	3	MAN	C4-C5-C6-O6
4	L	2	NAG	C3-C2-N2-C7
5	S	4	MAN	O5-C5-C6-O6

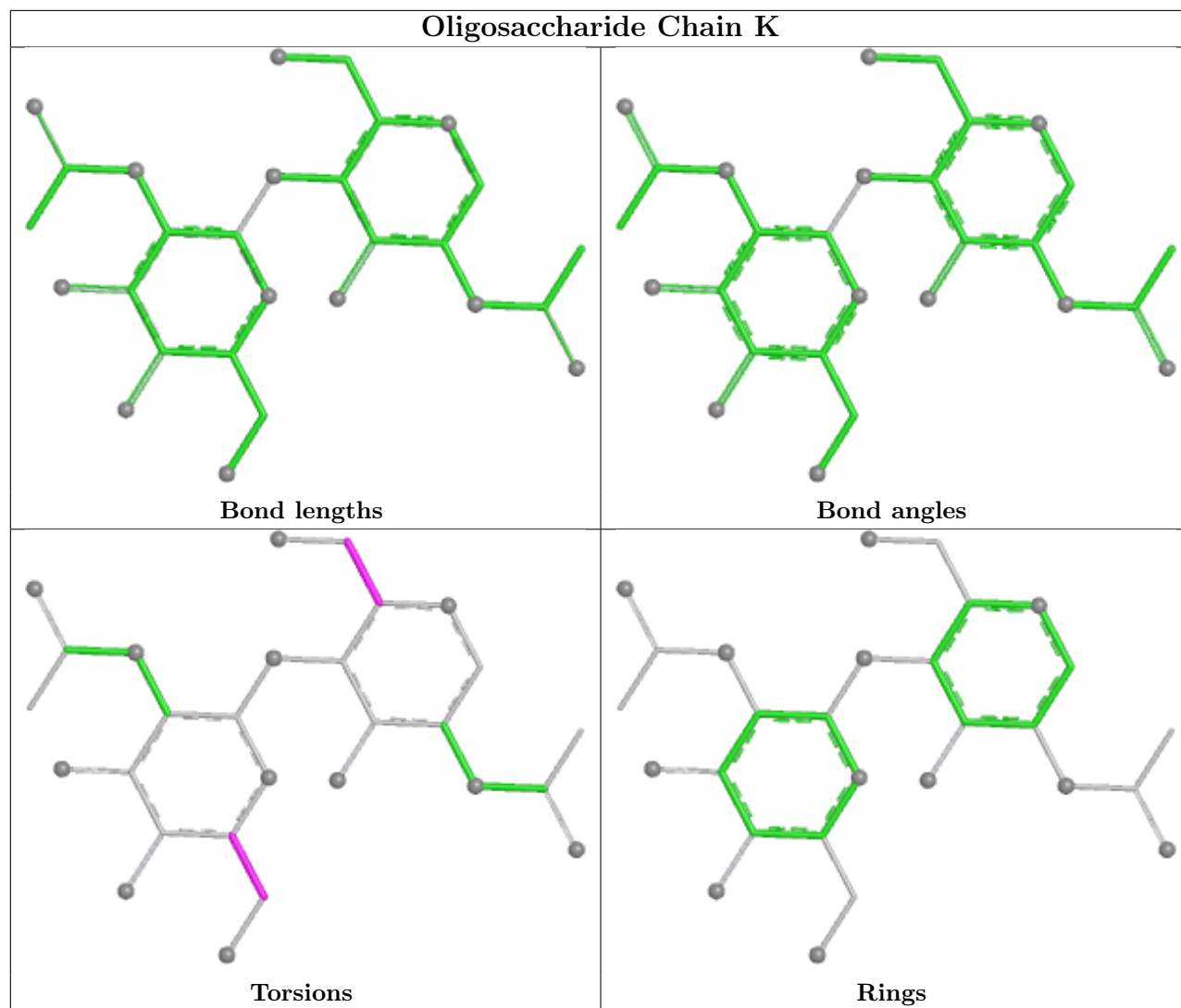
There are no ring outliers.

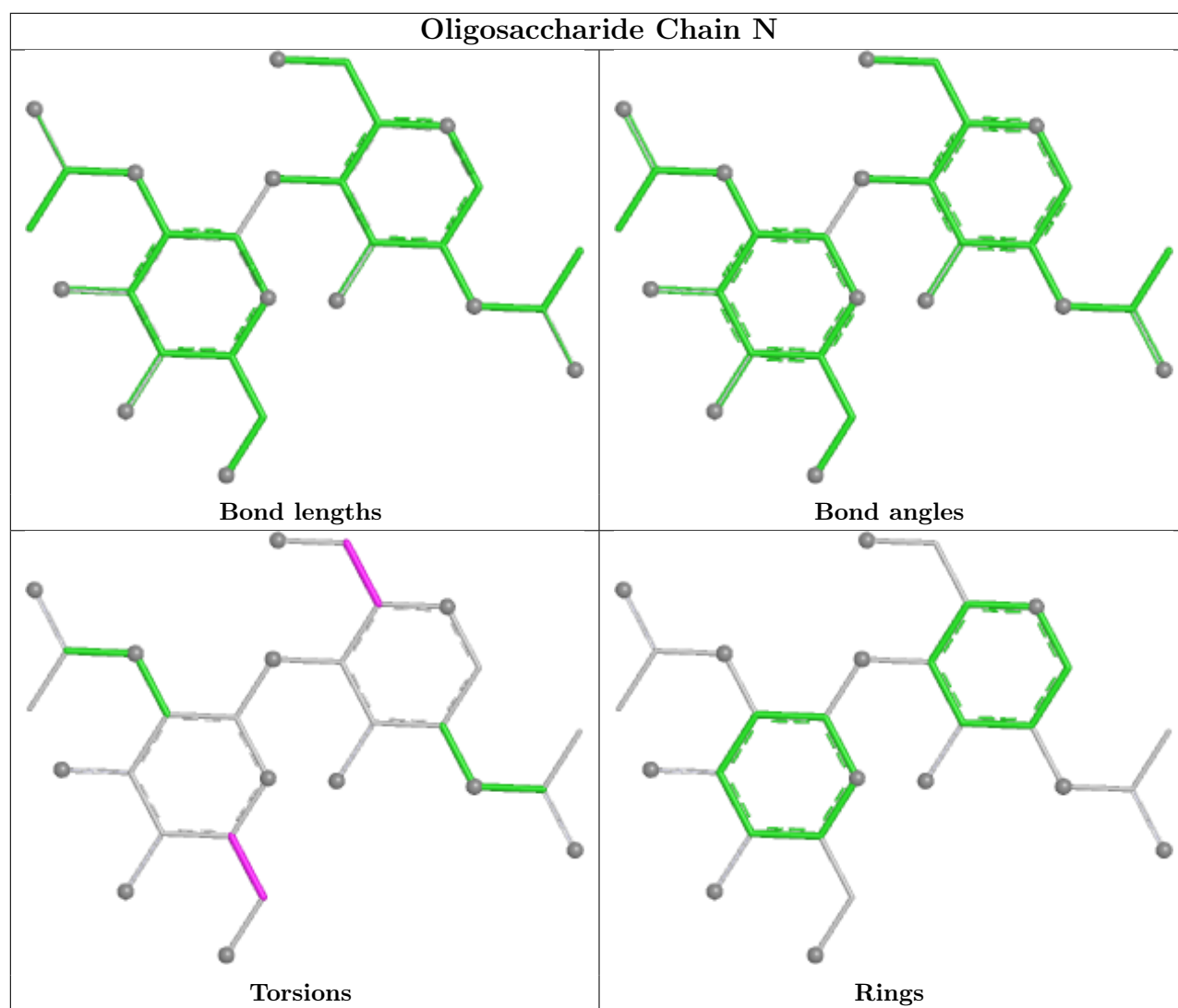
No monomer is involved in short contacts.

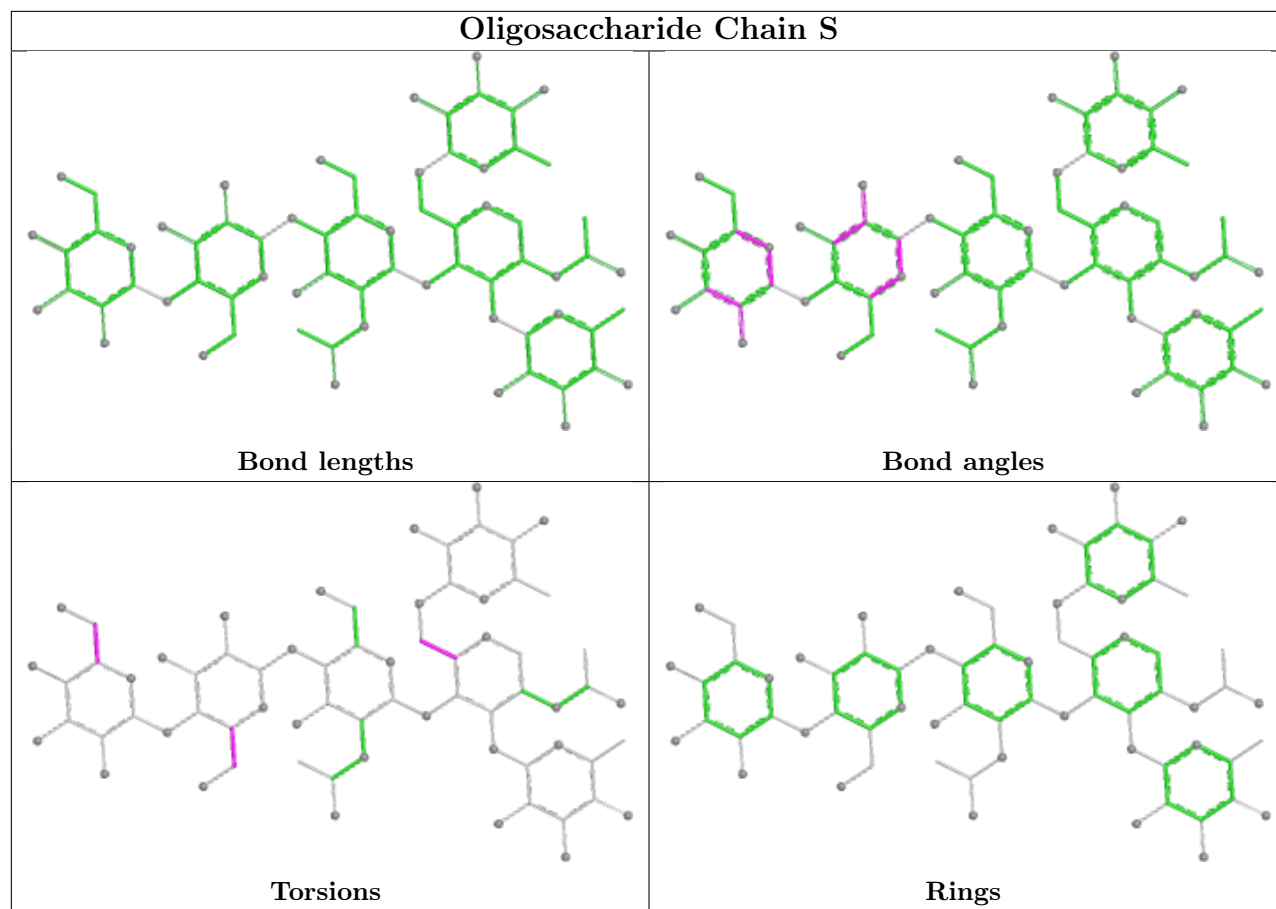
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for oligosaccharide.

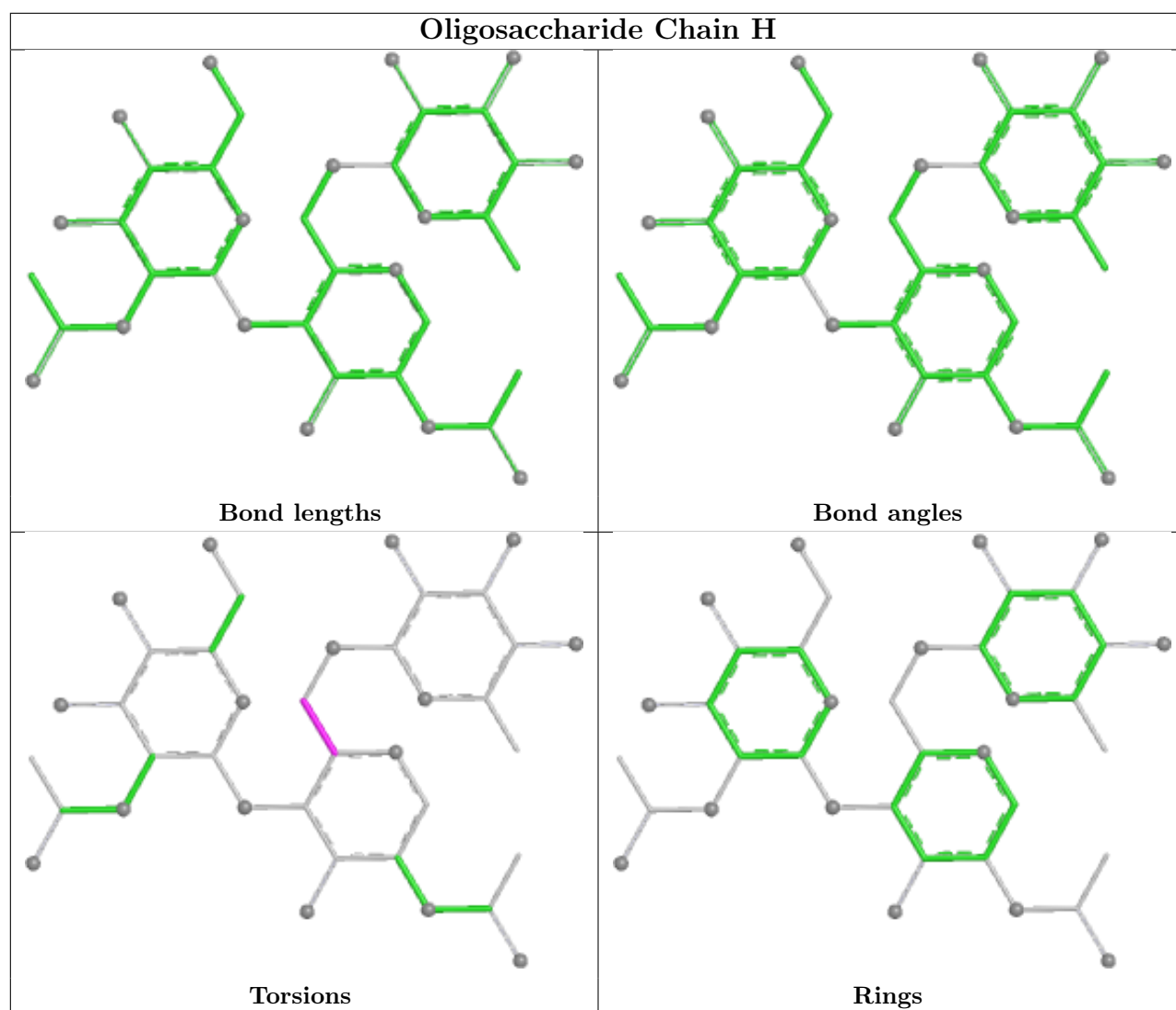












5.6 Ligand geometry [i](#)

12 ligands are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
7	NAG	B	704	2	14,14,15	0.23	0	17,19,21	0.45	0
7	NAG	B	701	2	14,14,15	0.23	0	17,19,21	0.44	0
7	NAG	B	702	2	14,14,15	0.27	0	17,19,21	0.44	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
7	NAG	B	705	2	14,14,15	0.25	0	17,19,21	0.47	0
7	NAG	B	703	2	14,14,15	0.23	0	17,19,21	0.43	0
7	NAG	C	201	3	14,14,15	0.28	0	17,19,21	0.34	0
7	NAG	E	704	2	14,14,15	0.24	0	17,19,21	0.43	0
7	NAG	E	701	2	14,14,15	0.25	0	17,19,21	0.45	0
7	NAG	E	705	2	14,14,15	0.22	0	17,19,21	0.35	0
7	NAG	E	703	2	14,14,15	0.24	0	17,19,21	0.44	0
7	NAG	F	201	3	14,14,15	0.28	0	17,19,21	0.45	0
7	NAG	E	702	2	14,14,15	0.20	0	17,19,21	0.42	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
7	NAG	B	704	2	-	2/6/23/26	0/1/1/1
7	NAG	B	701	2	-	1/6/23/26	0/1/1/1
7	NAG	B	702	2	-	0/6/23/26	0/1/1/1
7	NAG	B	705	2	-	0/6/23/26	0/1/1/1
7	NAG	B	703	2	-	1/6/23/26	0/1/1/1
7	NAG	C	201	3	-	2/6/23/26	0/1/1/1
7	NAG	E	704	2	-	2/6/23/26	0/1/1/1
7	NAG	E	701	2	-	0/6/23/26	0/1/1/1
7	NAG	E	705	2	-	0/6/23/26	0/1/1/1
7	NAG	E	703	2	-	2/6/23/26	0/1/1/1
7	NAG	F	201	3	-	0/6/23/26	0/1/1/1
7	NAG	E	702	2	-	0/6/23/26	0/1/1/1

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (10) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
7	C	201	NAG	C4-C5-C6-O6
7	C	201	NAG	O5-C5-C6-O6
7	E	703	NAG	O5-C5-C6-O6
7	E	703	NAG	C4-C5-C6-O6
7	E	704	NAG	O5-C5-C6-O6

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Mol	Chain	Res	Type	Atoms
7	B	704	NAG	C4-C5-C6-O6
7	B	704	NAG	O5-C5-C6-O6
7	B	703	NAG	O5-C5-C6-O6
7	E	704	NAG	C4-C5-C6-O6
7	B	701	NAG	O5-C5-C6-O6

There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
7	B	702	NAG	1	0
7	E	702	NAG	1	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ > 2		OWAB(Å ²)	Q < 0.9
1	A	170/185 (91%)	1.16	20 (11%)	9 6	38, 95, 137, 167	0
1	D	158/185 (85%)	1.41	38 (24%)	2 2	50, 105, 149, 167	0
2	B	644/677 (95%)	0.38	37 (5%)	29 19	14, 47, 110, 156	0
2	E	642/677 (94%)	0.46	44 (6%)	23 14	10, 45, 112, 161	0
3	C	114/157 (72%)	0.61	10 (8%)	15 10	20, 63, 116, 137	0
3	F	115/157 (73%)	0.82	12 (10%)	11 8	29, 71, 117, 141	0
All	All	1843/2038 (90%)	0.61	161 (8%)	16 11	10, 57, 127, 167	0

All (161) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	217	THR	5.7
2	E	698	TYR	5.5
2	E	217	THR	4.8
1	D	123	TYR	4.7
3	C	56	GLY	4.7
1	A	128	LEU	4.5
1	D	87	TYR	4.4
2	B	550	ASN	4.2
2	B	698	TYR	4.1
1	A	33	ASP	4.1
2	B	211	SER	3.9
2	E	550	ASN	3.9
1	D	109	PRO	3.8
1	D	76	ASP	3.8
1	A	204	LEU	3.8
2	E	677	ASN	3.7
1	D	101	THR	3.7
2	B	214	ALA	3.6
2	B	216	GLU	3.6

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Mol	Chain	Res	Type	RSRZ
2	B	215	ASN	3.6
3	F	131	ILE	3.6
2	B	212	LYS	3.6
2	E	658	ASN	3.6
2	B	630	PRO	3.6
1	A	123	TYR	3.5
2	B	115	CYS	3.5
1	A	48	TYR	3.5
2	E	574	VAL	3.4
2	B	131	ASN	3.4
1	D	91	ALA	3.4
1	D	186	GLY	3.4
1	D	189	VAL	3.3
2	E	407	GLY	3.3
1	A	115	CYS	3.2
2	E	37	ALA	3.2
1	A	189	VAL	3.2
2	B	650	LEU	3.1
3	F	115	VAL	3.1
2	E	122	GLN	3.1
2	B	114	THR	3.1
2	B	37	ALA	3.0
2	B	38	THR	3.0
1	D	84	ASN	3.0
2	E	191	GLY	3.0
2	B	113	TYR	3.0
2	B	408	ALA	3.0
1	D	41	LEU	3.0
1	A	87	TYR	3.0
2	B	122	GLN	3.0
2	E	40	LEU	2.9
2	E	610	SER	2.9
1	A	188	CYS	2.9
1	A	131	GLY	2.9
3	C	20	GLN	2.8
1	A	127	ASP	2.8
1	A	91	ALA	2.8
2	B	629	THR	2.8
1	A	30	VAL	2.8
2	E	627	ILE	2.8
1	A	138	LEU	2.8
2	E	216	GLU	2.7

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Mol	Chain	Res	Type	RSRZ
1	D	198	TYR	2.7
2	E	60	LEU	2.7
3	F	126	ILE	2.7
2	E	130	PHE	2.7
1	D	120	ASN	2.7
3	F	20	GLN	2.6
2	E	384	LYS	2.6
2	B	103	THR	2.6
1	A	146	ALA	2.6
1	D	74	SER	2.6
2	E	112	ARG	2.6
3	F	23	ALA	2.6
2	E	625	TYR	2.6
2	E	560	ILE	2.6
3	C	52	ASN	2.6
1	D	145	ILE	2.6
1	D	162	LYS	2.6
1	D	34	PHE	2.6
1	D	57	ASN	2.6
1	D	184	ASP	2.6
2	E	239	ASP	2.6
2	E	581	GLU	2.6
2	E	214	ALA	2.6
1	D	83	THR	2.6
1	D	163	LEU	2.5
1	D	88	ARG	2.5
2	E	585	ASN	2.5
2	E	628	SER	2.5
3	C	55	ASN	2.5
3	F	68	ASP	2.5
2	E	212	LYS	2.5
2	E	573	HIS	2.5
2	B	598	SER	2.5
2	B	70	THR	2.5
2	B	627	ILE	2.4
3	F	124	SER	2.4
2	E	629	THR	2.4
1	D	131	GLY	2.4
2	B	651	GLN	2.4
2	B	511	HIS	2.4
2	B	277	GLU	2.4
1	D	29	VAL	2.4

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Mol	Chain	Res	Type	RSRZ
2	B	512	PRO	2.4
3	C	66	THR	2.4
2	E	491	ASP	2.4
1	D	31	LEU	2.4
1	D	70	CYS	2.4
1	D	96	ILE	2.3
2	E	208	HIS	2.3
1	D	122	TYR	2.3
3	C	59	ILE	2.3
1	A	147	PRO	2.3
1	A	51	GLY	2.3
2	B	610	SER	2.3
3	F	22	VAL	2.3
1	D	181	ALA	2.3
1	A	107	SER	2.3
2	B	666	PHE	2.3
2	B	608	ASP	2.3
2	E	186	GLY	2.3
3	F	32	SER	2.3
2	B	669	LYS	2.2
2	E	121	SER	2.2
2	E	246	ASN	2.2
2	E	634	CYS	2.2
2	E	552	ILE	2.2
2	E	630	PRO	2.2
1	D	110	GLY	2.2
2	E	553	PRO	2.2
1	D	165	VAL	2.2
1	A	120	ASN	2.2
2	B	551	LEU	2.2
3	C	54	CYS	2.2
2	E	632	ARG	2.2
2	E	119	THR	2.2
1	D	48	TYR	2.2
1	A	129	ASP	2.2
2	B	91	LEU	2.1
1	D	50	LYS	2.1
2	B	581	GLU	2.1
2	B	658	ASN	2.1
1	D	78	ASP	2.1
1	D	86	VAL	2.1
2	E	68	VAL	2.1

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Mol	Chain	Res	Type	RSRZ
1	D	146	ALA	2.1
2	E	562	ALA	2.1
3	C	46	ILE	2.1
2	B	586	ALA	2.1
1	D	164	ASN	2.0
1	D	173	LEU	2.0
2	E	330	HIS	2.0
1	D	118	THR	2.0
3	F	27	CYS	2.0
3	F	57	VAL	2.0
2	E	107	SER	2.0
2	E	190	SER	2.0
2	B	451	ASP	2.0
3	C	116	GLY	2.0
3	F	25	PRO	2.0
3	C	19	ILE	2.0

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

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

6.3 Carbohydrates [i](#)

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

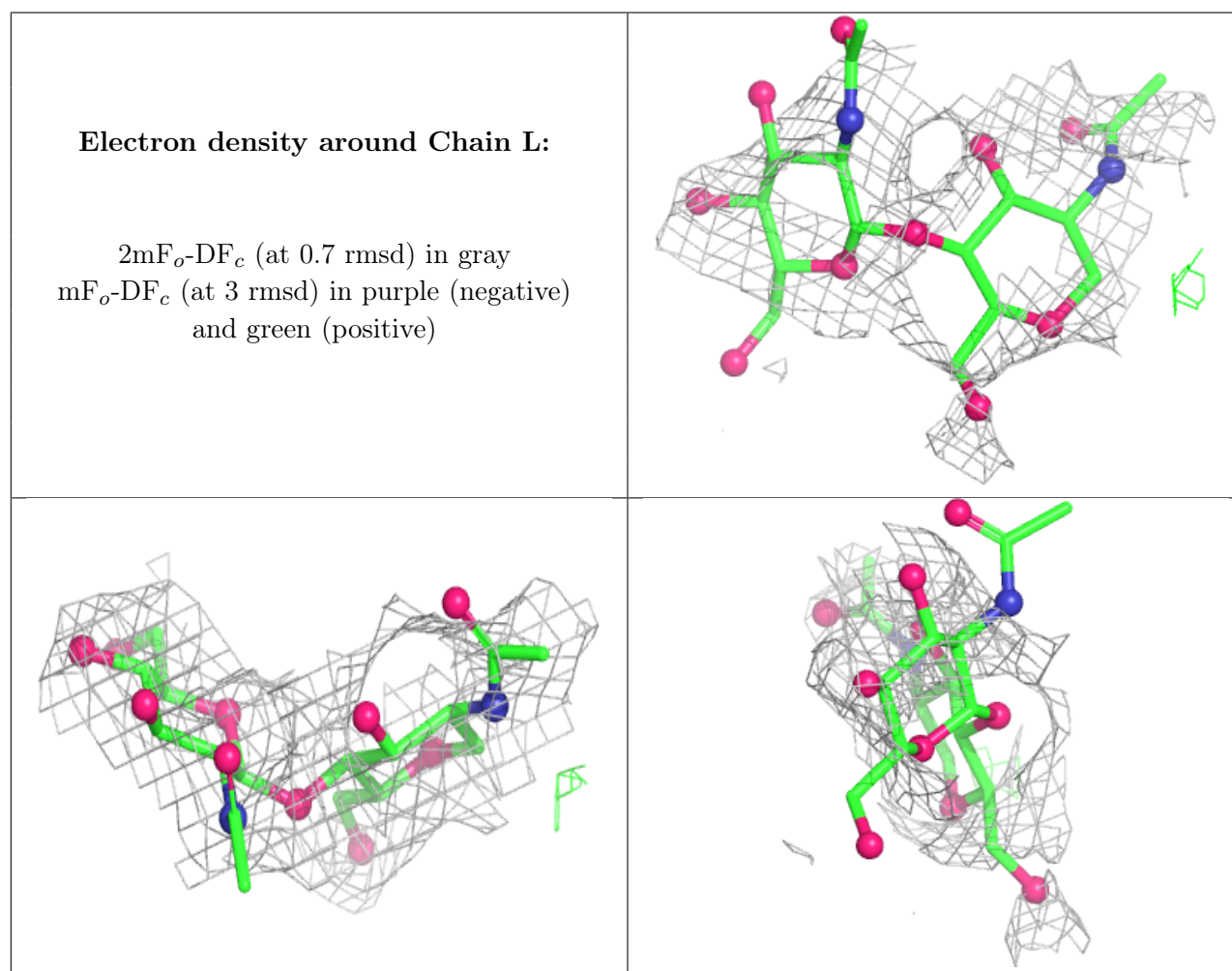
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
4	NAG	K	2	14/15	0.50	0.15	104,124,135,146	0
6	NAG	H	2	14/15	0.56	0.18	98,113,121,123	0
4	NAG	a	1	14/15	-	-	64,82,87,91	0
4	NAG	a	2	14/15	-	-	99,125,141,143	0
4	NAG	K	1	14/15	0.67	0.13	71,93,102,106	0
6	NAG	H	1	14/15	0.73	0.14	100,118,126,129	0
4	NAG	N	1	14/15	-	-	77,87,102,110	0
4	NAG	N	2	14/15	-	-	144,158,167,175	0
5	NAG	S	1	14/15	-	-	107,122,124,126	0
5	NAG	S	2	14/15	-	-	93,114,121,129	0
5	MAN	S	3	11/12	-	-	115,123,132,134	0
5	MAN	S	4	11/12	-	-	159,171,178,183	0

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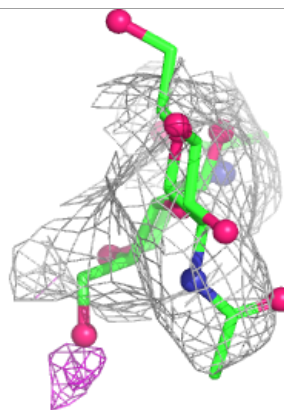
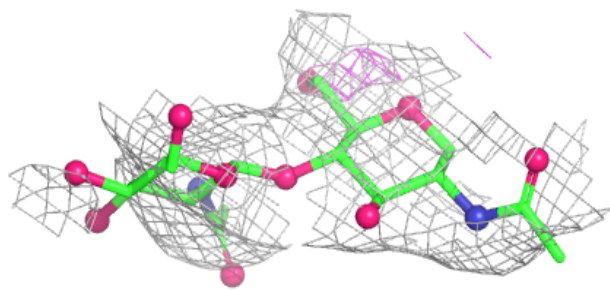
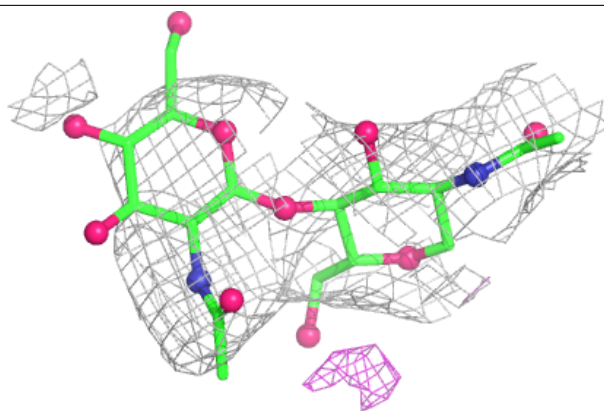
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
5	FUC	S	5	10/11	-	-	137,142,152,157	0
5	FUC	S	6	10/11	-	-	94,110,121,129	0
4	NAG	L	2	14/15	0.80	0.20	152,164,171,176	0
4	NAG	L	1	14/15	0.82	0.17	121,138,154,157	0
6	FUC	H	3	10/11	-	-	134,143,148,152	0

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.



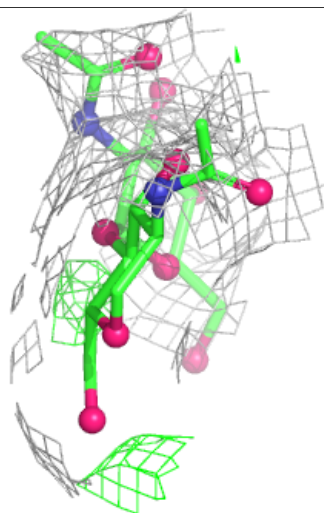
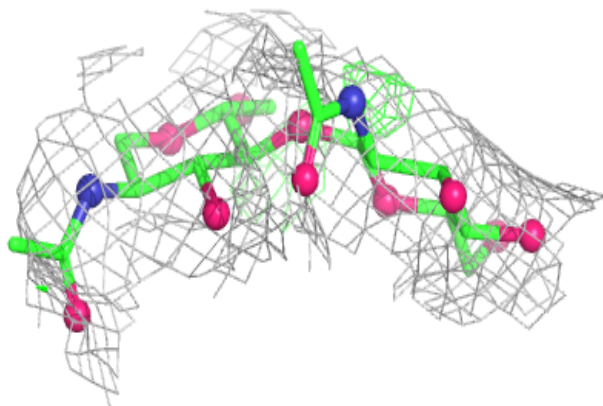
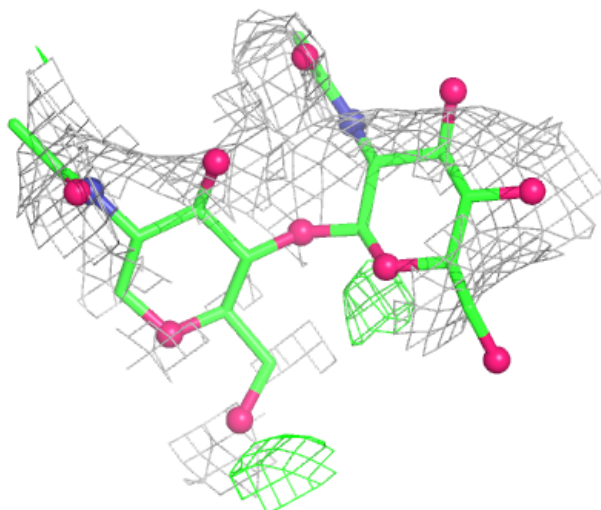
Electron density around Chain a:

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



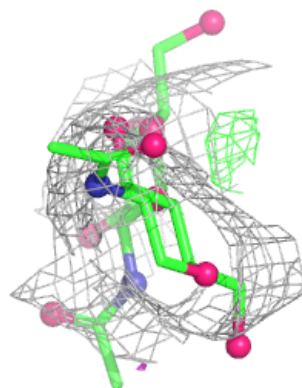
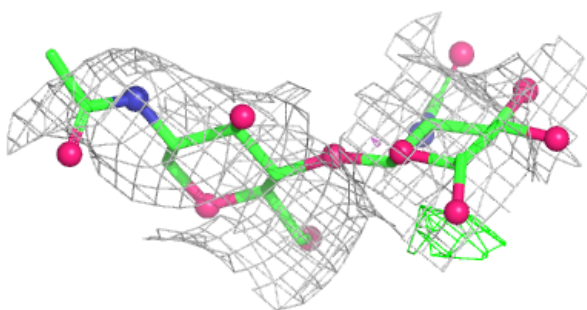
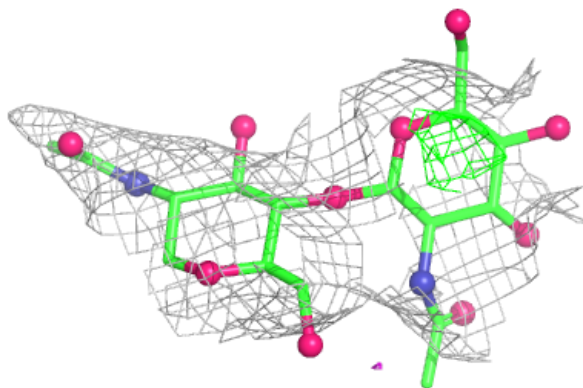
Electron density around Chain K:

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

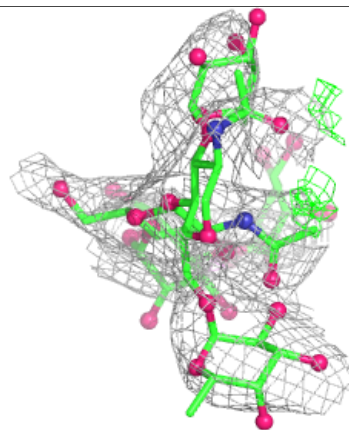
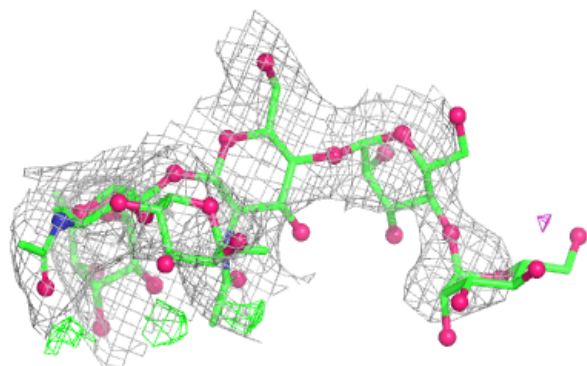
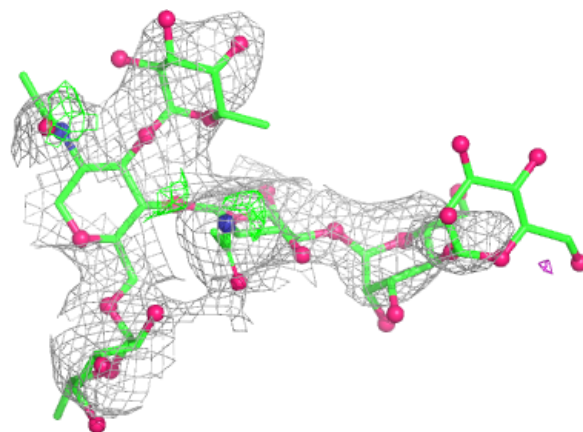


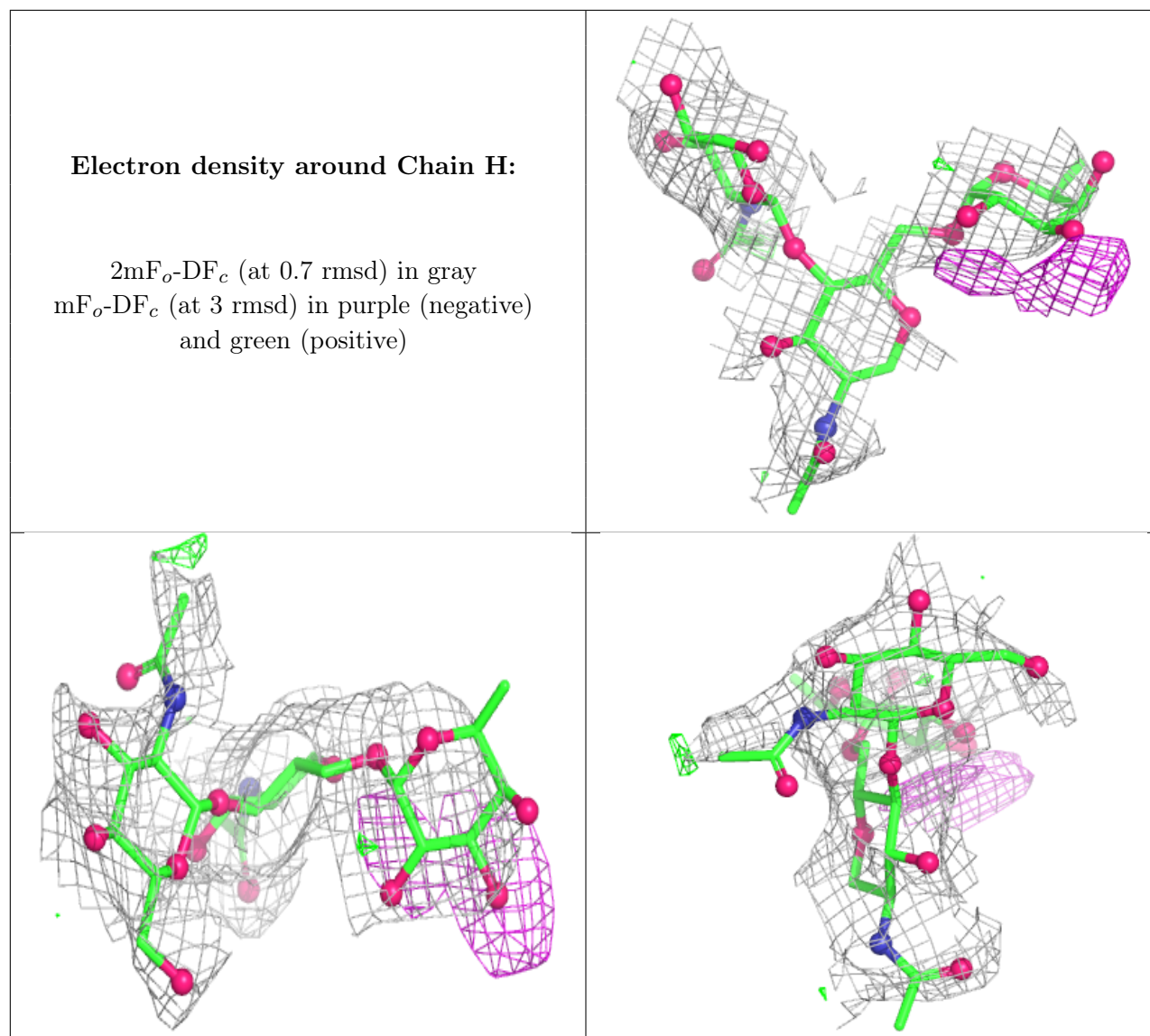
Electron density around Chain N:

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

**Electron density around Chain S:**

$2mF_o-DF_c$ (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
7	NAG	E	704	14/15	0.48	0.16	95,106,118,123	0
7	NAG	F	201	14/15	0.49	0.18	105,139,153,155	0
7	NAG	E	701	14/15	0.56	0.21	115,134,139,141	0
7	NAG	C	201	14/15	0.62	0.18	127,139,147,148	0
7	NAG	B	705	14/15	0.62	0.24	96,112,124,127	0
7	NAG	B	701	14/15	0.67	0.18	108,127,141,142	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
7	NAG	B	703	14/15	0.72	0.15	62,65,71,83	0
7	NAG	E	702	14/15	0.76	0.20	90,98,106,107	0
7	NAG	B	702	14/15	0.80	0.16	68,77,81,86	0
7	NAG	E	705	14/15	0.82	0.12	71,81,101,104	0
7	NAG	B	704	14/15	0.84	0.11	74,83,98,100	0
7	NAG	E	703	14/15	0.90	0.09	43,57,61,64	0

6.5 Other polymers [i](#)

There are no such residues in this entry.