



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 09:57 AM UTC

PDB ID : 6NM6 / pdb_00006nm6
Title : Crystal Structure of HIV-1 BG505 SOSIP.664 Prefusion Env Trimer Bound to N6 FR3-03 scFv in Complex with Crystallization Chaperones 3H109L Fab and 35O22 scFv at 3.2 Angstrom
Authors : Lai, Y.-T.; Kwong, P.D.
Deposited on : 2019-01-10
Resolution : 2.74 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

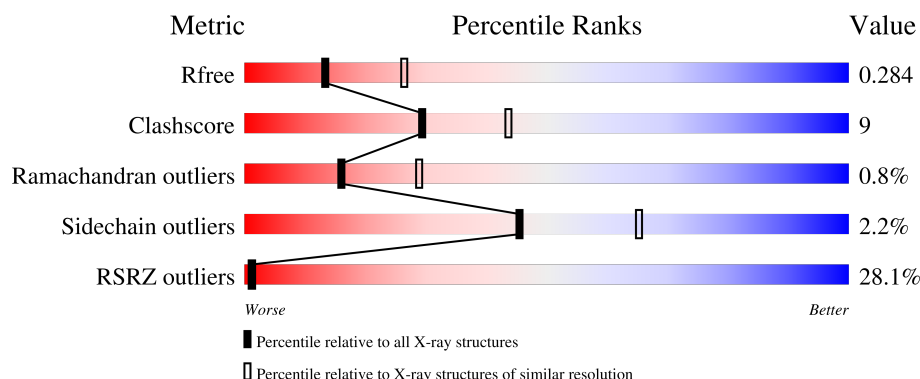
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.74 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	1819 (2.76-2.72)
Clashscore	190562	1866 (2.76-2.72)
Ramachandran outliers	187476	1830 (2.76-2.72)
Sidechain outliers	187428	1831 (2.76-2.72)
RSRZ outliers	180081	1819 (2.76-2.72)

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	B	153	
2	D	153	
3	E	130	
4	G	481	
5	H	244	

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Mol	Chain	Length	Quality of chain
6	L	217	 6% 74% 21% . .
7	U	145	 58% 72% 17% 12%
8	V	122	 74% 67% 13% . 19%
9	A	6	 33% 17% 50%
10	C	2	 100%
10	I	2	 100%
10	K	2	 100%
10	M	2	 100%
10	N	2	 100%
10	P	2	 100%
10	Q	2	 50% 50%
11	F	5	 60% 40%
12	J	4	 50% 50%
13	O	10	 10% 70% 20%

2 Entry composition

There are 15 unique types of molecules in this entry. The entry contains 11893 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 Envelope glycoprotein gp41.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	B	126	Total	C	N	O	S	0	0	0
			1001	635	172	188	6			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	559	PRO	ILE	engineered mutation	UNP Q2N0S6
B	605	CYS	THR	engineered mutation	UNP Q2N0S6

- Molecule 2 is a protein called 35O22 scFv heavy chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	D	128	Total	C	N	O	S	0	0	0
			994	628	169	192	5			

- Molecule 3 is a protein called 35O22 scFv light chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	E	110	Total	C	N	O	S	0	0	0
			836	525	138	167	6			

- Molecule 4 is a protein called Envelope glycoprotein gp120.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	G	429	Total	C	N	O	S	0	0	0
			3383	2131	598	627	27			

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
G	145	ALA	ASN	engineered mutation	UNP Q2N0S6
G	332	ASN	THR	engineered mutation	UNP Q2N0S6

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Chain	Residue	Modelled	Actual	Comment	Reference
G	501	CYS	ALA	engineered mutation	UNP Q2N0S6
G	509	ARG	-	expression tag	UNP Q2N0S6
G	510	ARG	-	expression tag	UNP Q2N0S6
G	511	ARG	-	expression tag	UNP Q2N0S6
G	512	ARG	-	expression tag	UNP Q2N0S6
G	513	ARG	-	expression tag	UNP Q2N0S6

- Molecule 5 is a protein called 3H109L Fab heavy chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
5	H	226	Total	C	N	O	S	0	0	0
			1715	1093	278	338	6			

- Molecule 6 is a protein called 3H109L Fab light chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
6	L	210	Total	C	N	O	S	0	0	0
			1598	1006	275	310	7			

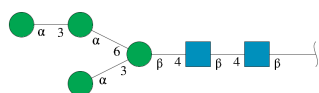
- Molecule 7 is a protein called N6 FR3-03 heavy chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
7	U	128	Total	C	N	O	S	0	0	0
			1005	637	180	184	4			

- Molecule 8 is a protein called N6 FR3-03 light chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
8	V	99	Total	C	N	O	S	0	0	0
			767	478	140	147	2			

- Molecule 9 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.



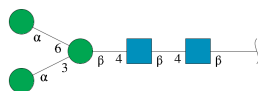
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
9	A	6	Total	C	N	O	0	0	0
			72	40	2	30			

- Molecule 10 is an oligosaccharide called 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.



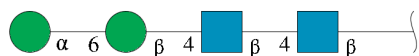
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
10	C	2	Total	C	N	O	0	0	0
			28	16	2	10			
10	I	2	Total	C	N	O	0	0	0
			28	16	2	10			
10	K	2	Total	C	N	O	0	0	0
			28	16	2	10			
10	M	2	Total	C	N	O	0	0	0
			28	16	2	10			
10	N	2	Total	C	N	O	0	0	0
			28	16	2	10			
10	P	2	Total	C	N	O	0	0	0
			28	16	2	10			
10	Q	2	Total	C	N	O	0	0	0
			28	16	2	10			

- Molecule 11 is an oligosaccharide called alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-6)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.



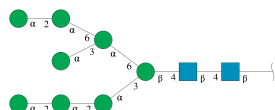
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
11	F	5	Total	C	N	O	0	0	0
			61	34	2	25			

- Molecule 12 is an oligosaccharide called alpha-D-mannopyranose-(1-6)-beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.



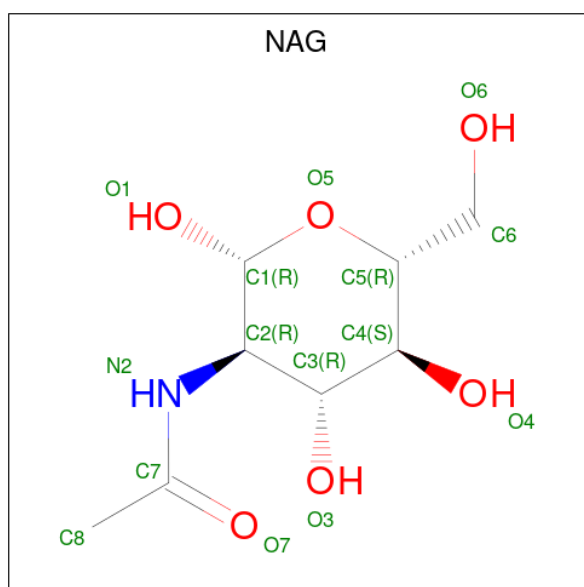
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
12	J	4	Total	C	N	O	0	0	0
			50	28	2	20			

- Molecule 13 is an oligosaccharide called alpha-D-mannopyranose-(1-2)-alpha-D-mannopyranose-(1-2)-alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-2)-alpha-D-mannopyranose-(1-6)-[alpha-D-mannopyranose-(1-3)]alpha-D-mannopyranose-(1-6)]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
13	O	10	Total	C	N	O	0	0	0
			116	64	2	50			

- Molecule 14 is 2-acetamido-2-deoxy-beta-D-glucopyranose (CCD ID: NAG) (formula: C₈H₁₅NO₆).



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

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

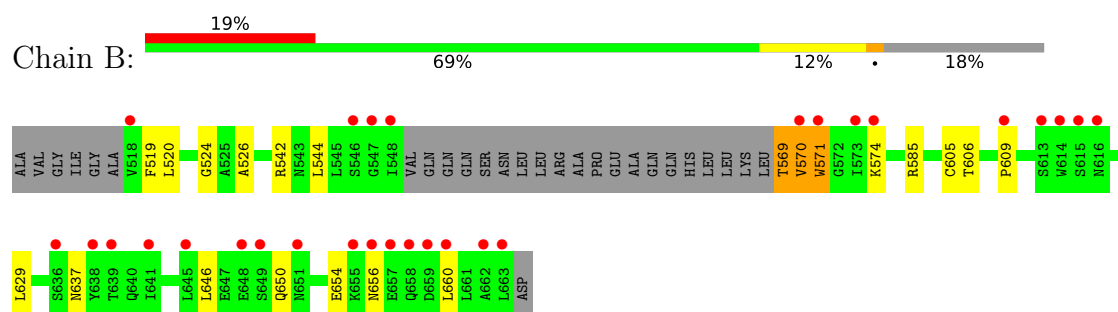
- Molecule 15 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
15	B	1	Total	O	0	0
			1	1		

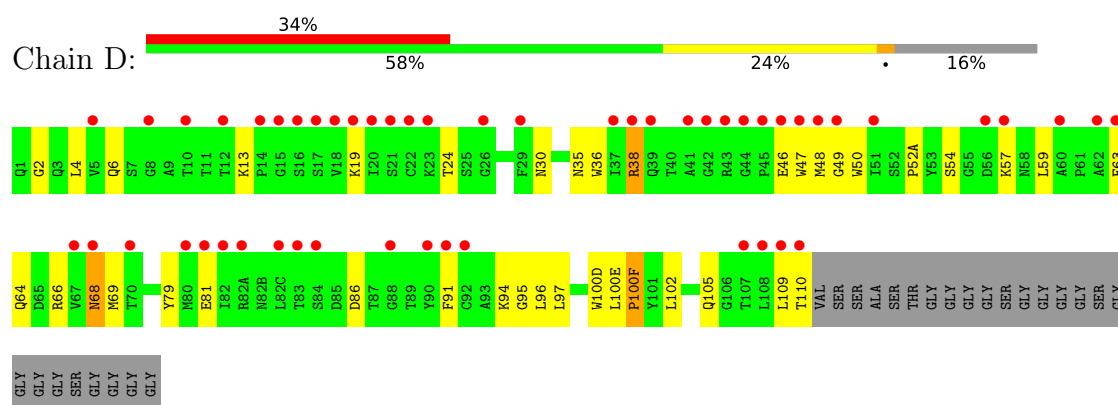
3 Residue-property plots [i](#)

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.

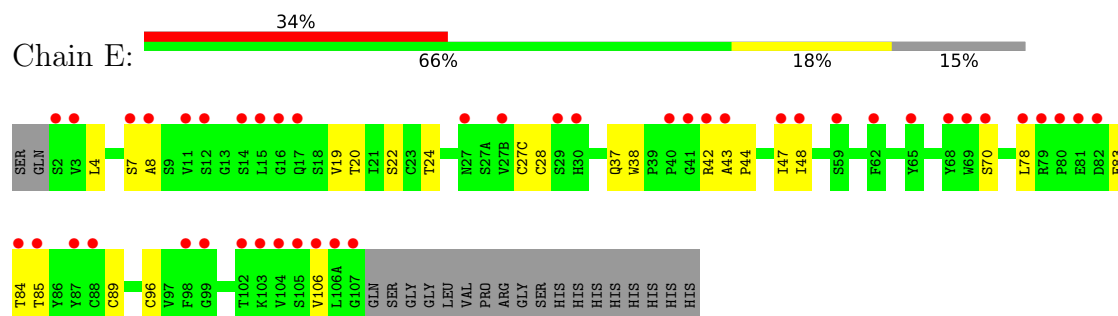
- Molecule 1: Envelope glycoprotein gp41



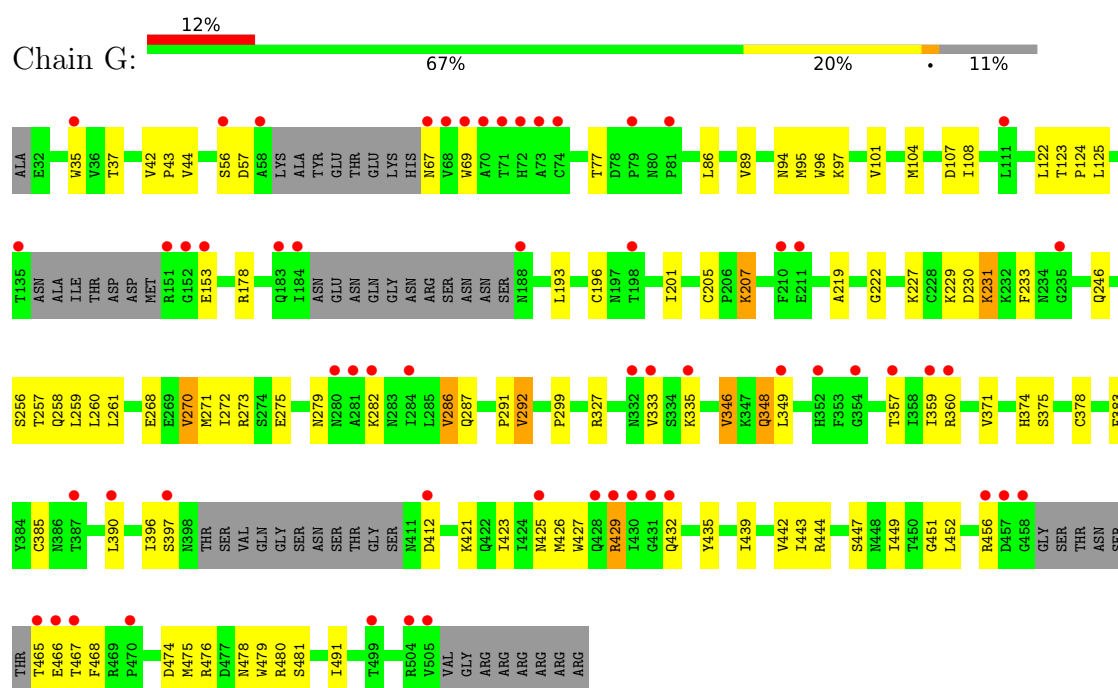
- Molecule 2: 35O22 scFv heavy chain



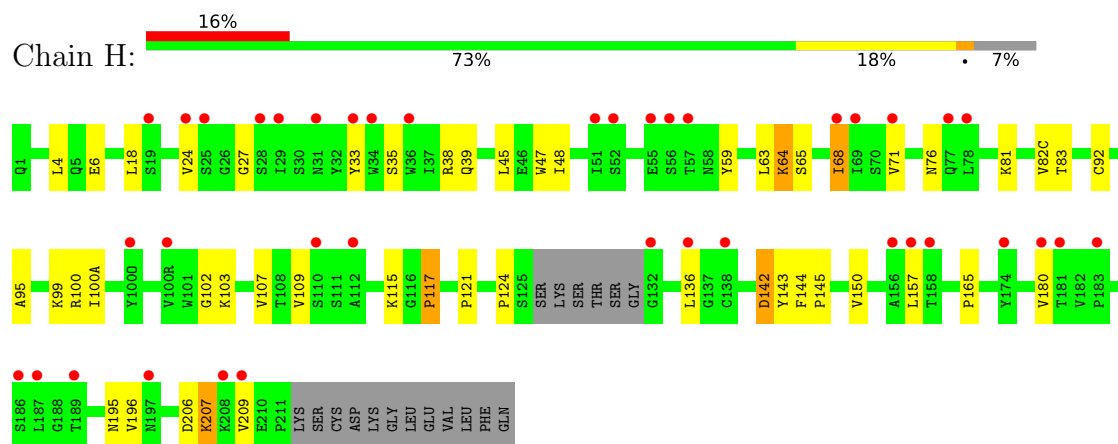
- Molecule 3: 35O22 scFv light chain



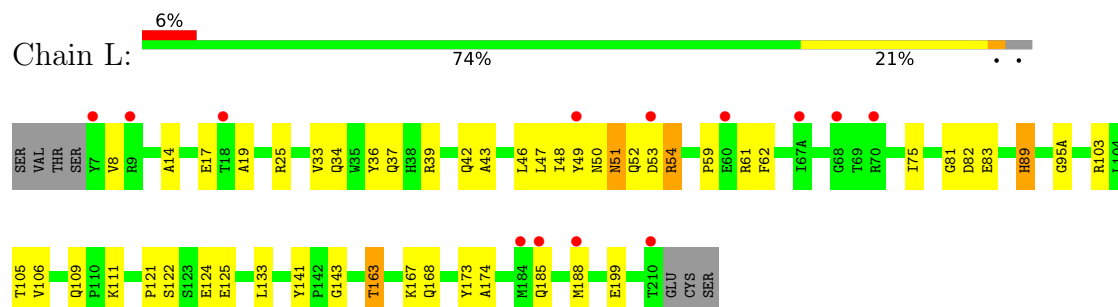
- Molecule 4: Envelope glycoprotein gp120



• Molecule 5: 3H109L Fab heavy chain

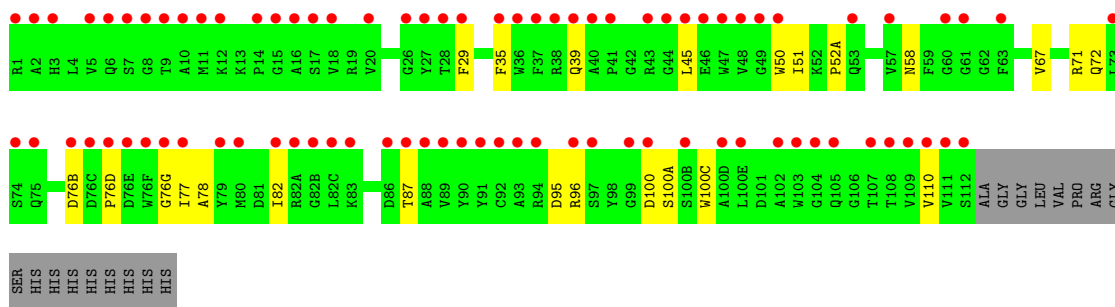


• Molecule 6: 3H109L Fab light chain

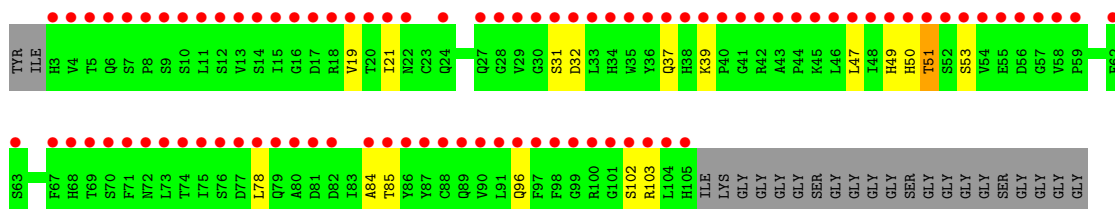
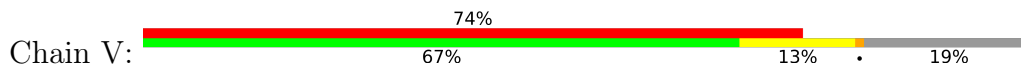


• Molecule 7: N6 FR3-03 heavy chain

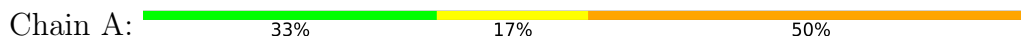




- Molecule 8: N6 FR3-03 light chain



- Molecule 9: 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



- Molecule 10: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose



- Molecule 10: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose



- Molecule 10: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose





- Molecule 10: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain M:  100%



- Molecule 10: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain N:  100%



- Molecule 10: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain P:  100%



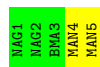
- Molecule 10: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain Q:  50% 50%

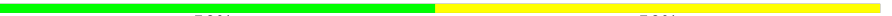


- Molecule 11: alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-6)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain F:  60% 40%



- Molecule 12: alpha-D-mannopyranose-(1-6)-beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain J:  50% 50%



- Molecule 13: α -D-mannopyranose-(1-2)- α -D-mannopyranose-(1-2)- α -D-mannopyranose-(1-3)-[α -D-mannopyranose-(1-2)- α -D-mannopyranose-(1-6)-[α -D-mannopyranose-(1-3)] α -D-mannopyranose-(1-6)] β -D-mannopyranose-(1-4)-2-acetamido-2-deoxy- β -D-glucopyranose-(1-4)-2-acetamido-2-deoxy- β -D-glucopyranose

Chain O:  10% 70% 20%

MAG1	MAG2	EMA3	MAN4	MAN5	MAN6	MAN7	MAN8	MAN9	MAN10
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4 Data and refinement statistics

Property	Value	Source
Space group	P 63	Depositor
Cell constants a, b, c, α , β , γ	128.24Å 128.24Å 315.77Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	37.02 – 2.74 37.02 – 2.74	Depositor EDS
% Data completeness (in resolution range)	45.1 (37.02-2.74) 45.1 (37.02-2.74)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.57 (at 2.72Å)	Xtriage
Refinement program	PHENIX 1.13_2998	Depositor
R, R_{free}	0.240 , 0.288 0.239 , 0.284	Depositor DCC
R_{free} test set	1735 reflections (2.25%)	wwPDB-VP
Wilson B-factor (Å ²)	33.8	Xtriage
Anisotropy	0.073	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.25 , 21.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.055 for h,-h-k,-l	Xtriage
F_o, F_c correlation	0.84	EDS
Total number of atoms	11893	wwPDB-VP
Average B, all atoms (Å ²)	57.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.26% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: BMA, NAG, 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	B	0.09	0/1019	0.25	0/1382
2	D	0.09	0/1021	0.30	0/1390
3	E	0.12	0/860	0.30	0/1175
4	G	0.10	0/3452	0.30	0/4683
5	H	0.10	0/1758	0.31	0/2397
6	L	0.10	0/1641	0.31	0/2239
7	U	0.09	0/1035	0.28	0/1408
8	V	0.09	0/787	0.30	0/1068
All	All	0.10	0/11573	0.30	0/15742

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	B	1001	0	982	16	0
2	D	994	0	952	25	0
3	E	836	0	785	13	0
4	G	3383	0	3334	71	0
5	H	1715	0	1685	30	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
6	L	1598	0	1548	29	0
7	U	1005	0	956	17	0
8	V	767	0	727	9	0
9	A	72	0	61	4	0
10	C	28	0	25	0	0
10	I	28	0	25	0	0
10	K	28	0	25	0	0
10	M	28	0	25	0	0
10	N	28	0	25	0	0
10	P	28	0	25	0	0
10	Q	28	0	25	1	0
11	F	61	0	52	0	0
12	J	50	0	43	2	0
13	O	116	0	97	2	0
14	B	42	0	39	0	0
14	D	14	0	13	1	0
14	G	42	0	39	0	0
15	B	1	0	0	0	0
All	All	11893	0	11488	201	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 9.

All (201) 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)
4:G:230:ASP:HB3	4:G:233:PHE:HB2	1.54	0.89
4:G:227:LYS:HD3	4:G:229:LYS:HZ1	1.52	0.75
4:G:201:ILE:HD11	4:G:435:TYR:HB2	1.70	0.74
5:H:59:TYR:HB2	5:H:64:LYS:HD2	1.68	0.74
1:B:605:CYS:HA	4:G:37:THR:HG22	1.72	0.71
1:B:520:LEU:HB3	1:B:524:GLY:HA3	1.72	0.71
4:G:335:LYS:HB3	4:G:412:ASP:HB3	1.72	0.71
4:G:456:ARG:O	7:U:58:ASN:ND2	2.24	0.69
1:B:585:ARG:NH2	4:G:491:ILE:O	2.25	0.69
3:E:78:LEU:HD13	3:E:106:VAL:HG23	1.75	0.68
4:G:257:THR:HG22	4:G:258:GLN:HG3	1.76	0.68
4:G:299:PRO:HG2	4:G:327:ARG:HB2	1.76	0.68
7:U:39:GLN:HB2	7:U:45:LEU:HD23	1.75	0.67
4:G:69:TRP:HE1	4:G:108:ILE:HD12	1.60	0.67
4:G:259:LEU:HD23	4:G:449:ILE:HG21	1.76	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:G:425:ASN:ND2	4:G:432:GLN:O	2.28	0.67
4:G:271:MET:HB3	4:G:273:ARG:HE	1.59	0.66
4:G:94:ASN:HB3	4:G:97:LYS:HG2	1.77	0.66
1:B:574:LYS:NZ	4:G:107:ASP:OD1	2.29	0.65
5:H:63:LEU:O	5:H:65:SER:N	2.29	0.64
4:G:282:LYS:NZ	7:U:100:ASP:OD1	2.23	0.63
5:H:157:LEU:HD21	5:H:180:VAL:HG11	1.79	0.63
6:L:54:ARG:NH2	6:L:62:PHE:O	2.30	0.63
1:B:571:TRP:HA	1:B:574:LYS:HB2	1.81	0.63
5:H:195:ASN:ND2	5:H:206:ASP:OD2	2.32	0.62
6:L:39:ARG:NH1	6:L:81:GLY:O	2.30	0.62
4:G:286:VAL:HG13	4:G:452:LEU:HB3	1.81	0.61
7:U:35:PHE:HE1	7:U:95:ASP:HB2	1.64	0.61
4:G:427:TRP:HZ3	4:G:475:MET:HG2	1.65	0.61
5:H:39:GLN:HB2	5:H:45:LEU:HD23	1.82	0.61
2:D:59:LEU:HD11	2:D:64:GLN:HA	1.82	0.61
2:D:36:TRP:HB3	2:D:48:MET:HE3	1.82	0.60
2:D:96:LEU:HG	2:D:97:LEU:HG	1.84	0.60
6:L:50:ASN:O	6:L:52:GLN:N	2.35	0.60
6:L:37:GLN:HB2	6:L:47:LEU:HD11	1.83	0.60
5:H:92:CYS:O	5:H:102:GLY:N	2.36	0.59
6:L:19:ALA:HB3	6:L:75:ILE:HB	1.84	0.59
4:G:292:VAL:HG13	4:G:449:ILE:HG13	1.85	0.58
2:D:19:LYS:HE3	2:D:79:TYR:HB3	1.85	0.57
9:A:2:NAG:H3	9:A:2:NAG:H83	1.86	0.57
4:G:475:MET:SD	4:G:478:ASN:ND2	2.78	0.57
7:U:51:ILE:HD11	7:U:71:ARG:HD2	1.85	0.57
7:U:72:GLN:HB3	7:U:77:ILE:HG23	1.86	0.57
5:H:100:ARG:NH2	13:O:4:MAN:O6	2.37	0.57
4:G:360:ARG:HB3	4:G:467:THR:HG22	1.87	0.57
4:G:456:ARG:NH1	4:G:466:GLU:OE2	2.38	0.57
4:G:279:ASN:OD1	7:U:100(C):TRP:NE1	2.37	0.56
3:E:37:GLN:HB2	3:E:47:ILE:HD11	1.85	0.56
5:H:136:LEU:HD13	5:H:209:VAL:HG21	1.87	0.56
2:D:38:ARG:HD2	2:D:46:GLU:HB3	1.88	0.56
4:G:258:GLN:NE2	4:G:371:VAL:O	2.34	0.56
5:H:35:SER:OG	5:H:47:TRP:NE1	2.39	0.56
4:G:260:LEU:HD12	4:G:451:GLY:HA3	1.88	0.55
2:D:57:LYS:HE3	2:D:69:MET:HE2	1.88	0.55
5:H:150:VAL:HG22	5:H:196:VAL:HG22	1.89	0.55
4:G:427:TRP:HD1	4:G:429:ARG:HD2	1.71	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:650:GLN:NE2	1:B:654:GLU:OE2	2.40	0.54
3:E:19:VAL:HG12	3:E:20:THR:H	1.71	0.54
3:E:47:ILE:HG22	3:E:48:ILE:HG13	1.89	0.54
2:D:35:ASN:ND2	2:D:100(D):TRP:O	2.40	0.54
6:L:8:VAL:HG11	6:L:103:ARG:HE	1.73	0.54
6:L:14:ALA:HB3	6:L:17:GLU:HG3	1.90	0.54
4:G:476:ARG:HA	4:G:479:TRP:CD1	2.43	0.54
5:H:117:PRO:HB3	5:H:143:TYR:HB3	1.89	0.53
6:L:59:PRO:HB2	6:L:61:ARG:HG2	1.89	0.53
6:L:83:GLU:OE1	6:L:167:LYS:NZ	2.41	0.53
8:V:39:LYS:HG2	8:V:84:ALA:HB2	1.91	0.53
8:V:37:GLN:HB2	8:V:47:LEU:HD11	1.89	0.53
1:B:571:TRP:H	1:B:571:TRP:CD1	2.27	0.53
6:L:124:GLU:N	6:L:124:GLU:OE1	2.40	0.53
2:D:68:ASN:HB2	2:D:81:GLU:HB2	1.90	0.53
4:G:42:VAL:HG23	4:G:44:VAL:HG12	1.90	0.53
4:G:346:VAL:HA	4:G:349:LEU:HD12	1.90	0.53
5:H:142:ASP:OD1	5:H:142:ASP:N	2.41	0.53
8:V:21:ILE:HD13	8:V:102:SER:HB2	1.91	0.53
4:G:439:ILE:HB	4:G:443:ILE:HD11	1.89	0.53
4:G:256:SER:OG	4:G:259:LEU:O	2.24	0.52
4:G:95:MET:HE3	4:G:480:ARG:HG2	1.91	0.52
5:H:18:LEU:HD11	5:H:107:VAL:HG11	1.92	0.52
4:G:69:TRP:NE1	4:G:108:ILE:HD12	2.24	0.52
1:B:519:PHE:HZ	1:B:542:ARG:HH22	1.57	0.52
4:G:374:HIS:HB3	4:G:385:CYS:HB2	1.92	0.52
4:G:256:SER:OG	4:G:257:THR:N	2.41	0.51
5:H:18:LEU:HB2	5:H:82(C):VAL:HG11	1.93	0.51
1:B:629:LEU:HD23	4:G:44:VAL:HG23	1.93	0.50
2:D:54:SER:HG	9:A:5:MAN:HO4	1.56	0.50
6:L:83:GLU:HG3	6:L:106:VAL:HG23	1.93	0.50
3:E:24:THR:HB	3:E:70:SER:HB3	1.94	0.50
8:V:19:VAL:HG21	8:V:78:LEU:HD13	1.92	0.50
2:D:66:ARG:NH1	2:D:86:ASP:OD2	2.45	0.50
5:H:68:ILE:HG23	5:H:81:LYS:HB2	1.93	0.50
5:H:99:LYS:HE2	5:H:100(A):ILE:HD11	1.94	0.50
6:L:34:GLN:HB2	6:L:89:HIS:HB3	1.93	0.50
2:D:100(E):LEU:HD12	2:D:100(F):PRO:HD2	1.95	0.49
3:E:37:GLN:HG3	3:E:84:THR:HG21	1.93	0.49
6:L:52:GLN:N	6:L:52:GLN:OE1	2.46	0.49
2:D:109:LEU:HG	2:D:110:THR:H	1.76	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:569:THR:O	1:B:569:THR:OG1	2.28	0.49
4:G:104:MET:O	4:G:108:ILE:HG12	2.13	0.49
4:G:257:THR:HB	4:G:375:SER:H	1.77	0.49
4:G:360:ARG:HH11	4:G:467:THR:HG21	1.78	0.49
13:O:1:NAG:H83	13:O:1:NAG:H3	1.94	0.49
4:G:478:ASN:O	4:G:481:SER:OG	2.30	0.48
5:H:165:PRO:HG2	6:L:163:THR:HG21	1.96	0.48
4:G:421:LYS:NZ	4:G:423:ILE:O	2.40	0.48
7:U:76(B):ASP:C	7:U:76(D):PRO:HD3	2.38	0.48
5:H:24:VAL:HG12	5:H:76:ASN:HB3	1.96	0.48
5:H:4:LEU:HG	5:H:24:VAL:HG23	1.96	0.48
6:L:185:GLN:HG2	6:L:188:MET:HE2	1.95	0.48
2:D:57:LYS:HE2	2:D:59:LEU:HD22	1.95	0.47
4:G:193:LEU:HB2	4:G:196:CYS:SG	2.54	0.47
6:L:52:GLN:NE2	6:L:53:ASP:OD1	2.47	0.47
2:D:57:LYS:HG3	2:D:69:MET:HE2	1.96	0.47
14:D:201:NAG:O7	14:D:201:NAG:O3	2.31	0.47
4:G:219:ALA:O	4:G:246:GLN:NE2	2.47	0.47
1:B:606:THR:HG21	1:B:646:LEU:HD22	1.96	0.47
1:B:609:PRO:HA	4:G:35:TRP:HA	1.96	0.47
3:E:83:GLU:HG3	3:E:106:VAL:HG12	1.96	0.47
4:G:123:THR:N	4:G:124:PRO:HD2	2.30	0.47
6:L:34:GLN:HG3	6:L:49:TYR:HA	1.95	0.47
5:H:33:TYR:HB2	5:H:95:ALA:O	2.15	0.47
1:B:526:ALA:HA	4:G:43:PRO:HB2	1.97	0.46
2:D:30:ASN:HA	2:D:52(A):PRO:HB2	1.96	0.46
7:U:95:ASP:OD1	7:U:96:ARG:N	2.48	0.46
3:E:7:SER:HB2	3:E:22:SER:H	1.80	0.46
6:L:47:LEU:O	6:L:48:ILE:HG13	2.16	0.46
8:V:85:THR:HG22	8:V:103:ARG:HB2	1.96	0.46
5:H:103:LYS:N	5:H:103:LYS:HD2	2.31	0.46
6:L:109:GLN:HB2	6:L:141:TYR:CE1	2.51	0.46
8:V:31:SER:O	8:V:51:THR:OG1	2.29	0.46
4:G:231:LYS:HB3	4:G:268:GLU:HG3	1.97	0.46
4:G:299:PRO:HA	4:G:442:VAL:HG13	1.97	0.46
4:G:333:VAL:HG21	4:G:390:LEU:HD21	1.98	0.45
4:G:101:VAL:HG13	4:G:479:TRP:HB2	1.98	0.45
4:G:378:CYS:HB3	4:G:383:PHE:CE1	2.52	0.45
4:G:205:CYS:HB3	4:G:207:LYS:HD2	1.99	0.45
6:L:83:GLU:HG2	6:L:105:THR:HA	1.99	0.45
4:G:122:LEU:HD13	4:G:125:LEU:HD12	1.99	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:H:121:PRO:HG3	5:H:207:LYS:HG2	1.97	0.44
2:D:2:GLY:O	2:D:102:LEU:HD21	2.18	0.44
4:G:357:THR:HG1	4:G:465:THR:N	2.15	0.44
2:D:48:MET:HG2	2:D:63:PHE:CD2	2.53	0.44
4:G:282:LYS:HZ3	7:U:100(A):SER:H	1.65	0.44
9:A:4:MAN:H3	9:A:5:MAN:H2	1.63	0.44
4:G:291:PRO:HG3	10:Q:1:NAG:O6	2.18	0.44
4:G:96:TRP:CG	4:G:275:GLU:HG2	2.53	0.44
3:E:38:TRP:CE2	3:E:44:PRO:HG3	2.53	0.44
5:H:103:LYS:HD2	5:H:103:LYS:H	1.83	0.43
7:U:29:PHE:CE2	7:U:52(A):PRO:HB3	2.53	0.43
7:U:100(C):TRP:CZ3	8:V:96:GLN:HB3	2.53	0.43
4:G:261:LEU:HD13	12:J:1:NAG:H82	1.99	0.43
4:G:56:SER:O	4:G:77:THR:N	2.35	0.43
5:H:144:PHE:HA	5:H:145:PRO:HA	1.79	0.43
5:H:38:ARG:HG2	5:H:48:ILE:HD11	2.01	0.43
5:H:83:THR:O	5:H:109:VAL:HG21	2.19	0.43
6:L:111:LYS:HD2	6:L:199:GLU:HG3	2.00	0.43
7:U:29:PHE:CD2	7:U:76(G):GLY:HA3	2.54	0.43
2:D:13:LYS:HD2	2:D:13:LYS:HA	1.68	0.43
2:D:47:TRP:HZ2	2:D:50:TRP:CD1	2.37	0.42
7:U:87:THR:HG23	7:U:110:VAL:HA	2.00	0.42
2:D:6:GLN:H	2:D:105:GLN:NE2	2.17	0.42
7:U:67:VAL:HG22	7:U:82:ILE:HG12	2.02	0.42
1:B:544:LEU:HD12	4:G:222:GLY:HA2	2.01	0.42
6:L:121:PRO:HD3	6:L:133:LEU:HD13	2.02	0.42
6:L:143:GLY:HA3	6:L:173:TYR:CG	2.54	0.42
4:G:86:LEU:HB3	4:G:89:VAL:HG21	2.02	0.42
2:D:47:TRP:CZ2	2:D:49:GLY:HA2	2.54	0.42
4:G:104:MET:SD	4:G:479:TRP:HB3	2.60	0.42
2:D:6:GLN:H	2:D:105:GLN:HE22	1.67	0.42
4:G:272:ILE:HG12	4:G:286:VAL:HB	2.01	0.42
8:V:32:ASP:HA	8:V:50:HIS:HA	2.02	0.42
5:H:24:VAL:CG1	5:H:76:ASN:HB3	2.50	0.42
4:G:153:GLU:O	4:G:178:ARG:HB2	2.20	0.41
5:H:6:GLU:OE1	5:H:6:GLU:N	2.51	0.41
2:D:94:LYS:HG2	2:D:95:GLY:O	2.19	0.41
4:G:272:ILE:HD12	4:G:348:GLN:HB3	2.02	0.41
6:L:168:GLN:OE1	6:L:174:ALA:HB2	2.20	0.41
4:G:261:LEU:HD23	4:G:449:ILE:HG22	2.02	0.41
6:L:122:SER:OG	6:L:125:GLU:HG2	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:E:84:THR:OG1	3:E:85:THR:N	2.53	0.41
4:G:359:ILE:HD12	4:G:468:PHE:HE1	1.84	0.41
2:D:4:LEU:HG	2:D:24:THR:HG22	2.02	0.41
2:D:50:TRP:CH2	9:A:4:MAN:H62	2.55	0.41
3:E:27(C):CYS:HA	3:E:28:CYS:HA	1.62	0.41
4:G:396:ILE:HG22	4:G:397:SER:N	2.35	0.41
8:V:49:HIS:O	8:V:53:SER:HB2	2.20	0.41
1:B:656:ASN:O	1:B:660:LEU:HG	2.21	0.41
4:G:282:LYS:HD3	4:G:282:LYS:HA	1.73	0.41
5:H:115:LYS:HA	5:H:115:LYS:HD2	1.85	0.41
7:U:35:PHE:CD2	7:U:50:TRP:HB3	2.55	0.41
1:B:637:ASN:OD1	1:B:637:ASN:N	2.54	0.41
6:L:33:VAL:HG12	6:L:51:ASN:OD1	2.21	0.41
4:G:447:SER:HB3	12:J:1:NAG:HN2	1.85	0.40
3:E:19:VAL:HG12	3:E:20:THR:N	2.36	0.40
5:H:27:GLY:O	5:H:76:ASN:ND2	2.54	0.40
6:L:42:GLN:HG2	6:L:43:ALA:H	1.86	0.40
4:G:270:VAL:HG23	4:G:287:GLN:O	2.21	0.40
6:L:36:TYR:HD1	6:L:46:LEU:HA	1.85	0.40
6:L:61:ARG:NH2	6:L:82:ASP:OD2	2.47	0.40
7:U:71:ARG:HA	7:U:78:ALA:HA	2.02	0.40
3:E:42:ARG:HB3	3:E:43:ALA:H	1.77	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	B	122/153 (80%)	113 (93%)	8 (7%)	1 (1%)	16	29
2	D	126/153 (82%)	113 (90%)	12 (10%)	1 (1%)	16	29
3	E	108/130 (83%)	88 (82%)	19 (18%)	1 (1%)	14	26

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
4	G	417/481 (87%)	378 (91%)	38 (9%)	1 (0%)	43	62
5	H	222/244 (91%)	198 (89%)	20 (9%)	4 (2%)	6	11
6	L	208/217 (96%)	189 (91%)	17 (8%)	2 (1%)	12	22
7	U	126/145 (87%)	119 (94%)	7 (6%)	0	100	100
8	V	97/122 (80%)	89 (92%)	7 (7%)	1 (1%)	12	22
All	All	1426/1645 (87%)	1287 (90%)	128 (9%)	11 (1%)	16	29

All (11) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
6	L	51	ASN
3	E	8	ALA
5	H	64	LYS
8	V	51	THR
5	H	117	PRO
4	G	474	ASP
1	B	570	VAL
5	H	124	PRO
6	L	95(A)	GLY
2	D	100(F)	PRO
5	H	68	ILE

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	B	108/129 (84%)	105 (97%)	3 (3%)	38	60
2	D	107/115 (93%)	104 (97%)	3 (3%)	38	60
3	E	96/113 (85%)	93 (97%)	3 (3%)	35	57
4	G	383/427 (90%)	371 (97%)	12 (3%)	35	57
5	H	196/212 (92%)	193 (98%)	3 (2%)	57	73
6	L	174/181 (96%)	170 (98%)	4 (2%)	44	65

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
7	U	102/115 (89%)	102 (100%)	0	100	100
8	V	87/94 (93%)	87 (100%)	0	100	100
All	All	1253/1386 (90%)	1225 (98%)	28 (2%)	45	65

All (28) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	569	THR
1	B	570	VAL
1	B	571	TRP
2	D	38	ARG
2	D	68	ASN
2	D	91	PHE
3	E	4	LEU
3	E	89	CYS
3	E	96	CYS
4	G	57	ASP
4	G	67	ASN
4	G	207	LYS
4	G	231	LYS
4	G	270	VAL
4	G	286	VAL
4	G	292	VAL
4	G	346	VAL
4	G	348	GLN
4	G	426	MET
4	G	429	ARG
4	G	444	ARG
5	H	71	VAL
5	H	142	ASP
5	H	207	LYS
6	L	25	ARG
6	L	54	ARG
6	L	89	HIS
6	L	163	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
4	G	85	HIS

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Mol	Chain	Res	Type
4	G	352	HIS
4	G	411	ASN
4	G	478	ASN
5	H	76	ASN
5	H	169	GLN
6	L	42	GLN
8	V	68	HIS

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 ⓘ

39 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	$\# Z > 2$
9	NAG	A	1	9,4	14,14,15	0.30	0	17,19,21	0.47	0
9	NAG	A	2	9	14,14,15	0.46	0	17,19,21	1.35	2 (11%)
9	BMA	A	3	9	11,11,12	0.67	0	15,15,17	0.85	0
9	MAN	A	4	9	11,11,12	0.93	1 (9%)	15,15,17	1.21	2 (13%)
9	MAN	A	5	9	11,11,12	1.47	1 (9%)	15,15,17	1.67	3 (20%)
9	MAN	A	6	9	11,11,12	0.59	0	15,15,17	1.03	2 (13%)
10	NAG	C	1	10,4	14,14,15	0.19	0	17,19,21	0.47	0
10	NAG	C	2	10	14,14,15	0.24	0	17,19,21	0.47	0
11	NAG	F	1	11,4	14,14,15	0.26	0	17,19,21	0.43	0
11	NAG	F	2	11	14,14,15	0.24	0	17,19,21	0.39	0
11	BMA	F	3	11	11,11,12	0.67	0	15,15,17	0.82	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
11	MAN	F	4	11	11,11,12	0.77	1 (9%)	15,15,17	1.08	2 (13%)
11	MAN	F	5	11	11,11,12	0.70	0	15,15,17	1.02	2 (13%)
10	NAG	I	1	10,4	14,14,15	0.28	0	17,19,21	0.50	0
10	NAG	I	2	10	14,14,15	0.27	0	17,19,21	0.38	0
12	NAG	J	1	12,4	14,14,15	0.30	0	17,19,21	0.55	0
12	NAG	J	2	12	14,14,15	0.25	0	17,19,21	0.47	0
12	BMA	J	3	12	11,11,12	0.63	0	15,15,17	0.72	0
12	MAN	J	4	12	11,11,12	0.65	0	15,15,17	1.10	2 (13%)
10	NAG	K	1	10,4	14,14,15	0.25	0	17,19,21	0.40	0
10	NAG	K	2	10	14,14,15	0.28	0	17,19,21	0.37	0
10	NAG	M	1	10,4	14,14,15	0.23	0	17,19,21	0.46	0
10	NAG	M	2	10	14,14,15	0.26	0	17,19,21	0.42	0
10	NAG	N	1	10,4	14,14,15	0.29	0	17,19,21	0.45	0
10	NAG	N	2	10	14,14,15	0.25	0	17,19,21	0.51	0
13	NAG	O	1	13,4	14,14,15	0.37	0	17,19,21	1.50	2 (11%)
13	MAN	O	10	13	11,11,12	1.06	2 (18%)	15,15,17	1.44	3 (20%)
13	NAG	O	2	13	14,14,15	0.22	0	17,19,21	0.42	0
13	BMA	O	3	13	11,11,12	0.96	1 (9%)	15,15,17	0.92	0
13	MAN	O	4	13	11,11,12	0.74	1 (9%)	15,15,17	1.34	2 (13%)
13	MAN	O	5	13	11,11,12	0.65	0	15,15,17	0.92	1 (6%)
13	MAN	O	6	13	11,11,12	0.70	0	15,15,17	0.83	1 (6%)
13	MAN	O	7	13	11,11,12	0.70	0	15,15,17	0.99	2 (13%)
13	MAN	O	8	13	11,11,12	0.69	0	15,15,17	1.04	1 (6%)
13	MAN	O	9	13	11,11,12	0.76	1 (9%)	15,15,17	1.27	2 (13%)
10	NAG	P	1	10,4	14,14,15	0.28	0	17,19,21	0.50	0
10	NAG	P	2	10	14,14,15	0.28	0	17,19,21	0.52	0
10	NAG	Q	1	10,4	14,14,15	0.95	1 (7%)	17,19,21	1.68	1 (5%)
10	NAG	Q	2	10	14,14,15	0.27	0	17,19,21	0.46	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
9	NAG	A	1	9,4	-	0/6/23/26	0/1/1/1
9	NAG	A	2	9	-	6/6/23/26	0/1/1/1
9	BMA	A	3	9	-	2/2/19/22	0/1/1/1
9	MAN	A	4	9	-	1/2/19/22	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
9	MAN	A	5	9	-	2/2/19/22	0/1/1/1
9	MAN	A	6	9	-	0/2/19/22	0/1/1/1
10	NAG	C	1	10,4	-	0/6/23/26	0/1/1/1
10	NAG	C	2	10	-	2/6/23/26	0/1/1/1
11	NAG	F	1	11,4	-	2/6/23/26	0/1/1/1
11	NAG	F	2	11	-	0/6/23/26	0/1/1/1
11	BMA	F	3	11	-	2/2/19/22	0/1/1/1
11	MAN	F	4	11	-	2/2/19/22	0/1/1/1
11	MAN	F	5	11	-	2/2/19/22	0/1/1/1
10	NAG	I	1	10,4	-	2/6/23/26	0/1/1/1
10	NAG	I	2	10	-	3/6/23/26	0/1/1/1
12	NAG	J	1	12,4	-	2/6/23/26	0/1/1/1
12	NAG	J	2	12	-	2/6/23/26	0/1/1/1
12	BMA	J	3	12	-	2/2/19/22	0/1/1/1
12	MAN	J	4	12	-	2/2/19/22	0/1/1/1
10	NAG	K	1	10,4	-	2/6/23/26	0/1/1/1
10	NAG	K	2	10	-	2/6/23/26	0/1/1/1
10	NAG	M	1	10,4	-	2/6/23/26	0/1/1/1
10	NAG	M	2	10	-	0/6/23/26	0/1/1/1
10	NAG	N	1	10,4	-	0/6/23/26	0/1/1/1
10	NAG	N	2	10	-	2/6/23/26	0/1/1/1
13	NAG	O	1	13,4	-	6/6/23/26	0/1/1/1
13	MAN	O	10	13	-	0/2/19/22	0/1/1/1
13	NAG	O	2	13	-	2/6/23/26	0/1/1/1
13	BMA	O	3	13	-	0/2/19/22	0/1/1/1
13	MAN	O	4	13	-	2/2/19/22	0/1/1/1
13	MAN	O	5	13	-	0/2/19/22	0/1/1/1
13	MAN	O	6	13	-	0/2/19/22	0/1/1/1
13	MAN	O	7	13	-	0/2/19/22	0/1/1/1
13	MAN	O	8	13	-	0/2/19/22	0/1/1/1
13	MAN	O	9	13	-	1/2/19/22	0/1/1/1
10	NAG	P	1	10,4	-	2/6/23/26	0/1/1/1
10	NAG	P	2	10	-	0/6/23/26	0/1/1/1
10	NAG	Q	1	10,4	-	2/6/23/26	0/1/1/1
10	NAG	Q	2	10	-	2/6/23/26	0/1/1/1

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
9	A	5	MAN	C1-C2	4.13	1.62	1.52
10	Q	1	NAG	O5-C1	3.39	1.49	1.43
13	O	9	MAN	C1-C2	2.25	1.57	1.52
13	O	10	MAN	C1-C2	2.25	1.57	1.52
13	O	10	MAN	C2-C3	2.18	1.55	1.52
13	O	4	MAN	C1-C2	2.17	1.57	1.52
11	F	4	MAN	C1-C2	2.12	1.57	1.52
13	O	3	BMA	O5-C1	-2.08	1.40	1.43
9	A	4	MAN	O5-C1	-2.01	1.40	1.43

All (30) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
10	Q	1	NAG	C1-O5-C5	6.51	120.91	112.19
13	O	1	NAG	C2-N2-C7	4.80	129.34	122.90
9	A	2	NAG	C2-N2-C7	4.55	129.00	122.90
9	A	5	MAN	C1-O5-C5	4.05	117.62	112.19
13	O	4	MAN	C1-O5-C5	3.89	117.39	112.19
13	O	9	MAN	C1-O5-C5	3.67	117.10	112.19
9	A	5	MAN	C1-C2-C3	3.56	114.83	109.64
13	O	8	MAN	C1-O5-C5	3.07	116.30	112.19
12	J	4	MAN	C1-O5-C5	3.06	116.29	112.19
13	O	10	MAN	C1-C2-C3	3.04	114.08	109.64
9	A	6	MAN	C1-O5-C5	2.87	116.03	112.19
13	O	1	NAG	C1-C2-N2	2.67	114.64	110.43
11	F	5	MAN	C1-O5-C5	2.47	115.49	112.19
11	F	4	MAN	C1-O5-C5	2.45	115.46	112.19
13	O	10	MAN	C1-O5-C5	2.44	115.46	112.19
13	O	7	MAN	C1-O5-C5	2.41	115.42	112.19
9	A	4	MAN	C3-C4-C5	2.37	114.53	110.23
9	A	4	MAN	O2-C2-C3	-2.26	105.48	110.15
13	O	4	MAN	O2-C2-C3	-2.24	105.50	110.15
13	O	10	MAN	O2-C2-C3	-2.20	105.59	110.15
9	A	6	MAN	O2-C2-C3	-2.19	105.61	110.15
13	O	9	MAN	O2-C2-C3	-2.17	105.65	110.15
12	J	4	MAN	O2-C2-C3	-2.16	105.68	110.15
13	O	7	MAN	O2-C2-C3	-2.15	105.70	110.15
9	A	5	MAN	O2-C2-C3	-2.14	105.71	110.15
11	F	5	MAN	O2-C2-C3	-2.14	105.72	110.15
9	A	2	NAG	C1-C2-N2	2.13	113.79	110.43
13	O	5	MAN	O2-C2-C3	-2.08	105.84	110.15
11	F	4	MAN	O2-C2-C3	-2.06	105.89	110.15
13	O	6	MAN	O2-C2-C3	-2.04	105.94	110.15

There are no chirality outliers.

All (59) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
10	I	2	NAG	O5-C5-C6-O6
10	M	1	NAG	O5-C5-C6-O6
9	A	2	NAG	O5-C5-C6-O6
12	J	1	NAG	O5-C5-C6-O6
10	Q	2	NAG	O5-C5-C6-O6
11	F	4	MAN	O5-C5-C6-O6
9	A	5	MAN	O5-C5-C6-O6
12	J	2	NAG	O5-C5-C6-O6
9	A	3	BMA	O5-C5-C6-O6
10	I	1	NAG	O5-C5-C6-O6
9	A	3	BMA	C4-C5-C6-O6
10	I	2	NAG	C4-C5-C6-O6
12	J	4	MAN	O5-C5-C6-O6
11	F	3	BMA	C4-C5-C6-O6
12	J	2	NAG	C4-C5-C6-O6
13	O	1	NAG	O5-C5-C6-O6
13	O	1	NAG	C4-C5-C6-O6
10	M	1	NAG	C4-C5-C6-O6
9	A	2	NAG	C4-C5-C6-O6
11	F	4	MAN	C4-C5-C6-O6
12	J	1	NAG	C4-C5-C6-O6
13	O	4	MAN	O5-C5-C6-O6
10	Q	2	NAG	C4-C5-C6-O6
12	J	4	MAN	C4-C5-C6-O6
9	A	2	NAG	C8-C7-N2-C2
9	A	2	NAG	O7-C7-N2-C2
10	Q	1	NAG	C8-C7-N2-C2
10	Q	1	NAG	O7-C7-N2-C2
13	O	1	NAG	C8-C7-N2-C2
13	O	1	NAG	O7-C7-N2-C2
9	A	5	MAN	C4-C5-C6-O6
11	F	5	MAN	C4-C5-C6-O6
13	O	4	MAN	C4-C5-C6-O6
10	P	1	NAG	O5-C5-C6-O6
11	F	3	BMA	O5-C5-C6-O6
11	F	5	MAN	O5-C5-C6-O6
12	J	3	BMA	O5-C5-C6-O6
10	K	1	NAG	O5-C5-C6-O6
10	K	1	NAG	C4-C5-C6-O6
11	F	1	NAG	C4-C5-C6-O6

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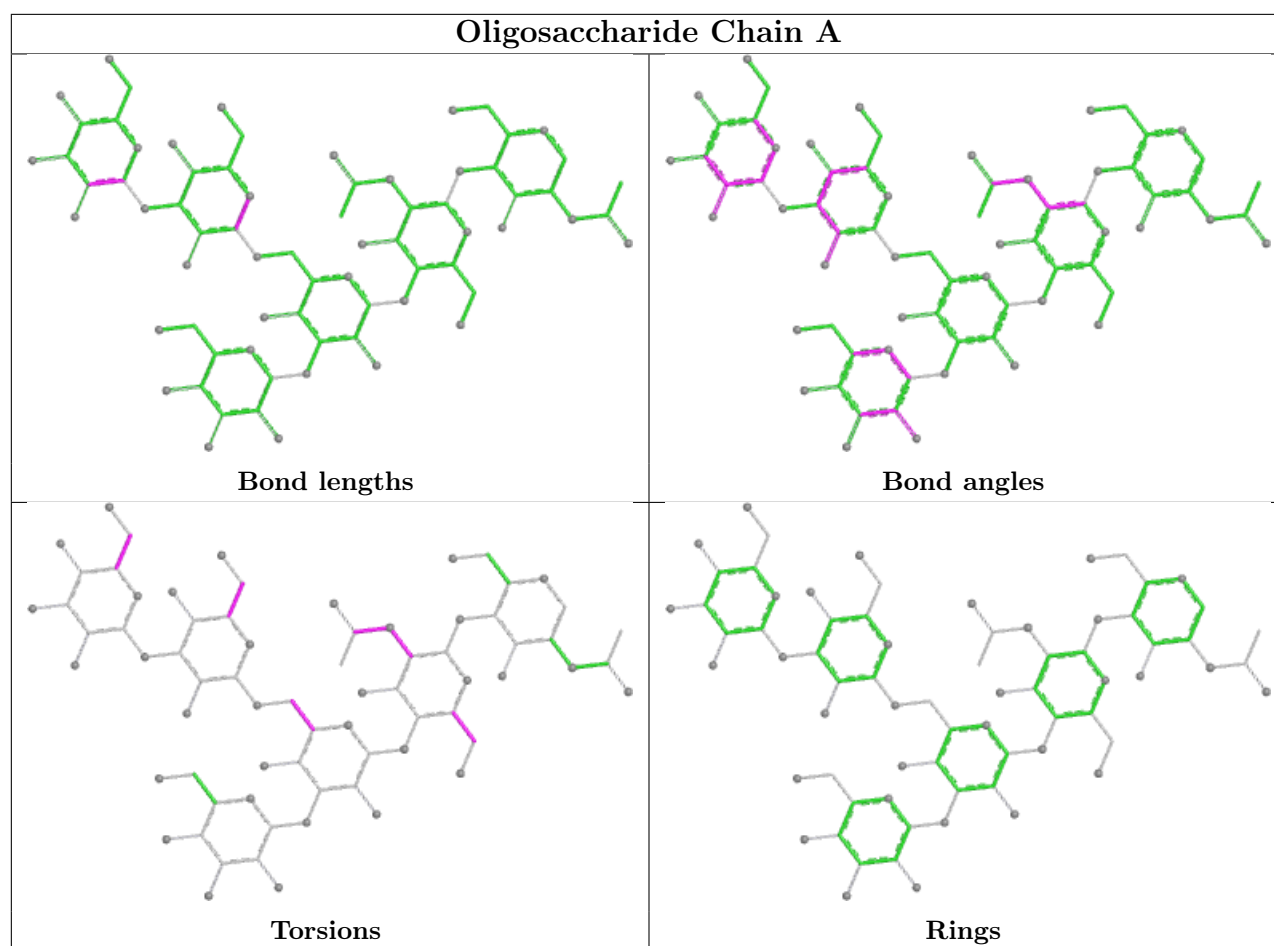
Mol	Chain	Res	Type	Atoms
10	I	1	NAG	C4-C5-C6-O6
13	O	2	NAG	C4-C5-C6-O6
11	F	1	NAG	O5-C5-C6-O6
10	C	2	NAG	C4-C5-C6-O6
10	K	2	NAG	C4-C5-C6-O6
13	O	2	NAG	O5-C5-C6-O6
10	N	2	NAG	C4-C5-C6-O6
10	P	1	NAG	C4-C5-C6-O6
10	K	2	NAG	O5-C5-C6-O6
10	C	2	NAG	O5-C5-C6-O6
13	O	9	MAN	C4-C5-C6-O6
12	J	3	BMA	C4-C5-C6-O6
10	N	2	NAG	O5-C5-C6-O6
9	A	2	NAG	C1-C2-N2-C7
10	I	2	NAG	C1-C2-N2-C7
13	O	1	NAG	C1-C2-N2-C7
9	A	2	NAG	C3-C2-N2-C7
13	O	1	NAG	C3-C2-N2-C7
9	A	4	MAN	C4-C5-C6-O6

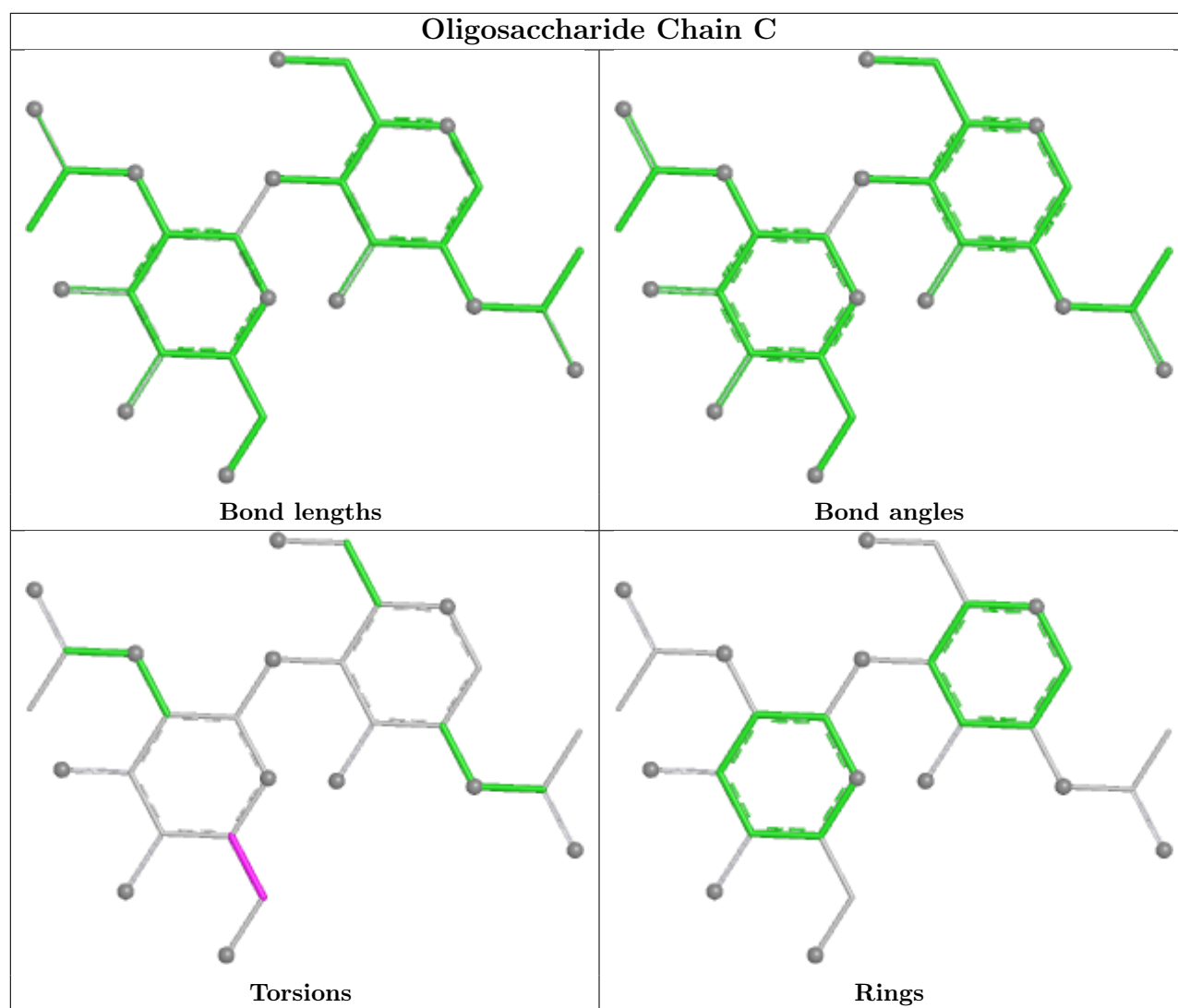
There are no ring outliers.

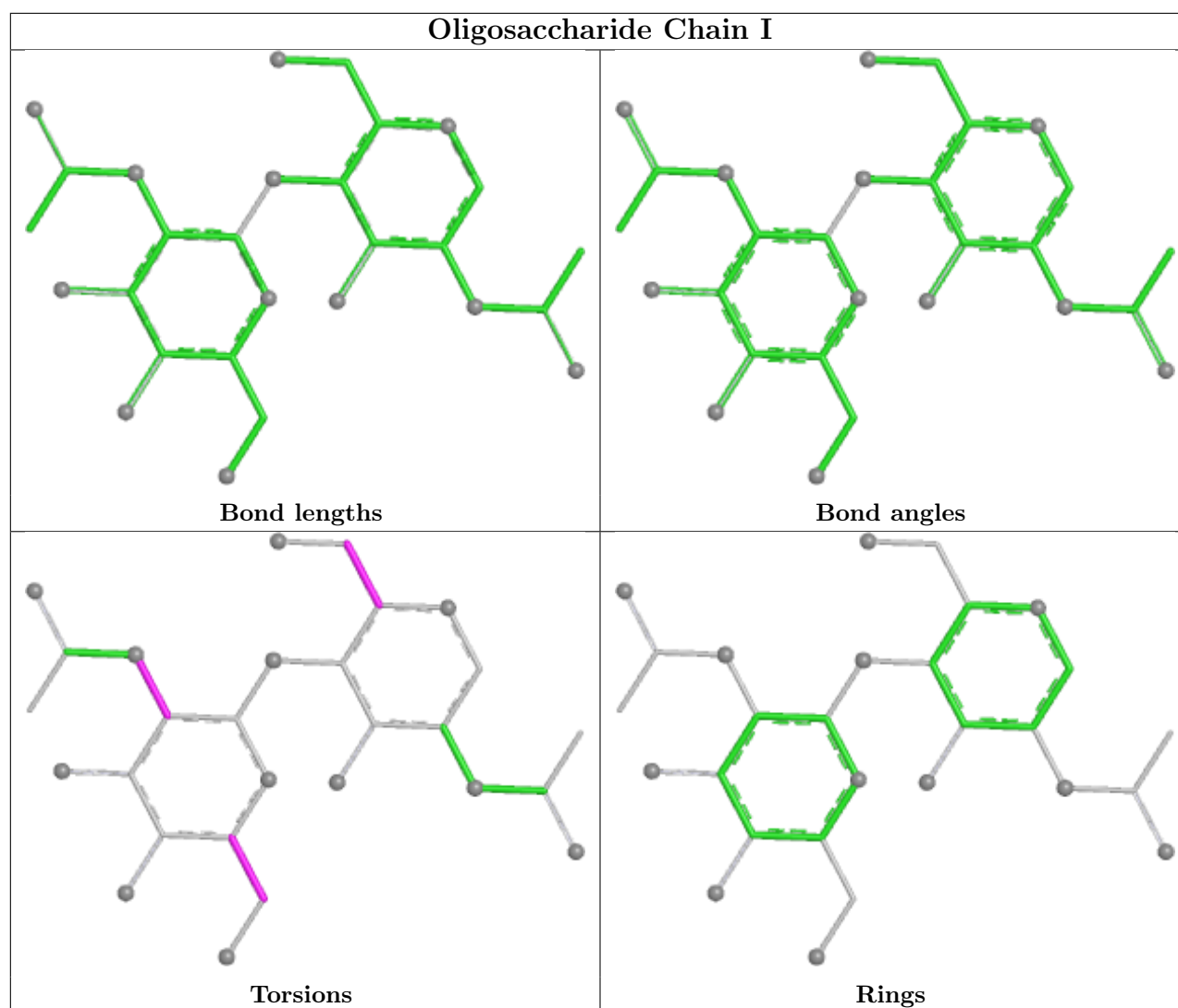
7 monomers are involved in 9 short contacts:

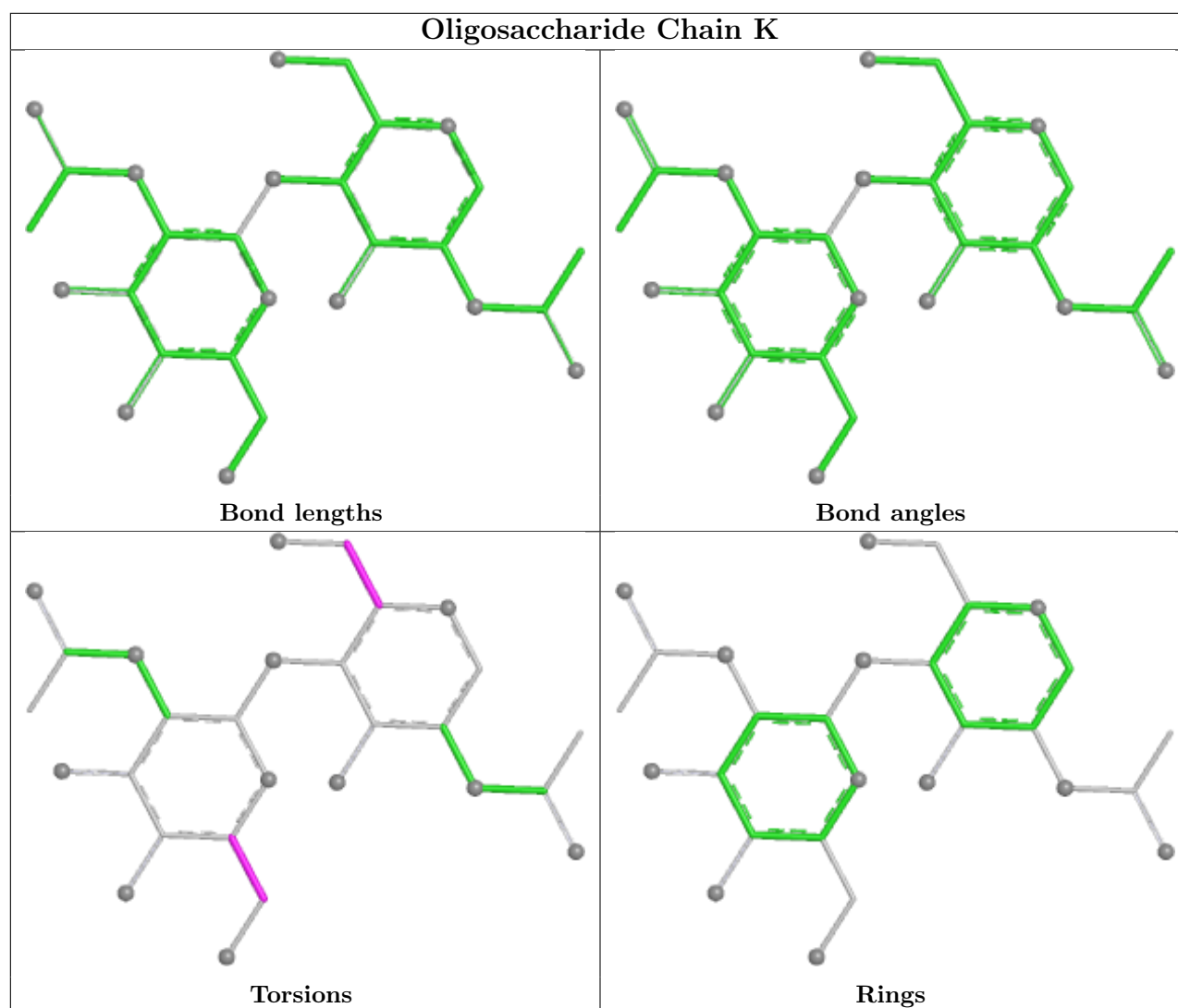
Mol	Chain	Res	Type	Clashes	Symm-Clashes
9	A	5	MAN	2	0
12	J	1	NAG	2	0
9	A	4	MAN	2	0
13	O	1	NAG	1	0
9	A	2	NAG	1	0
13	O	4	MAN	1	0
10	Q	1	NAG	1	0

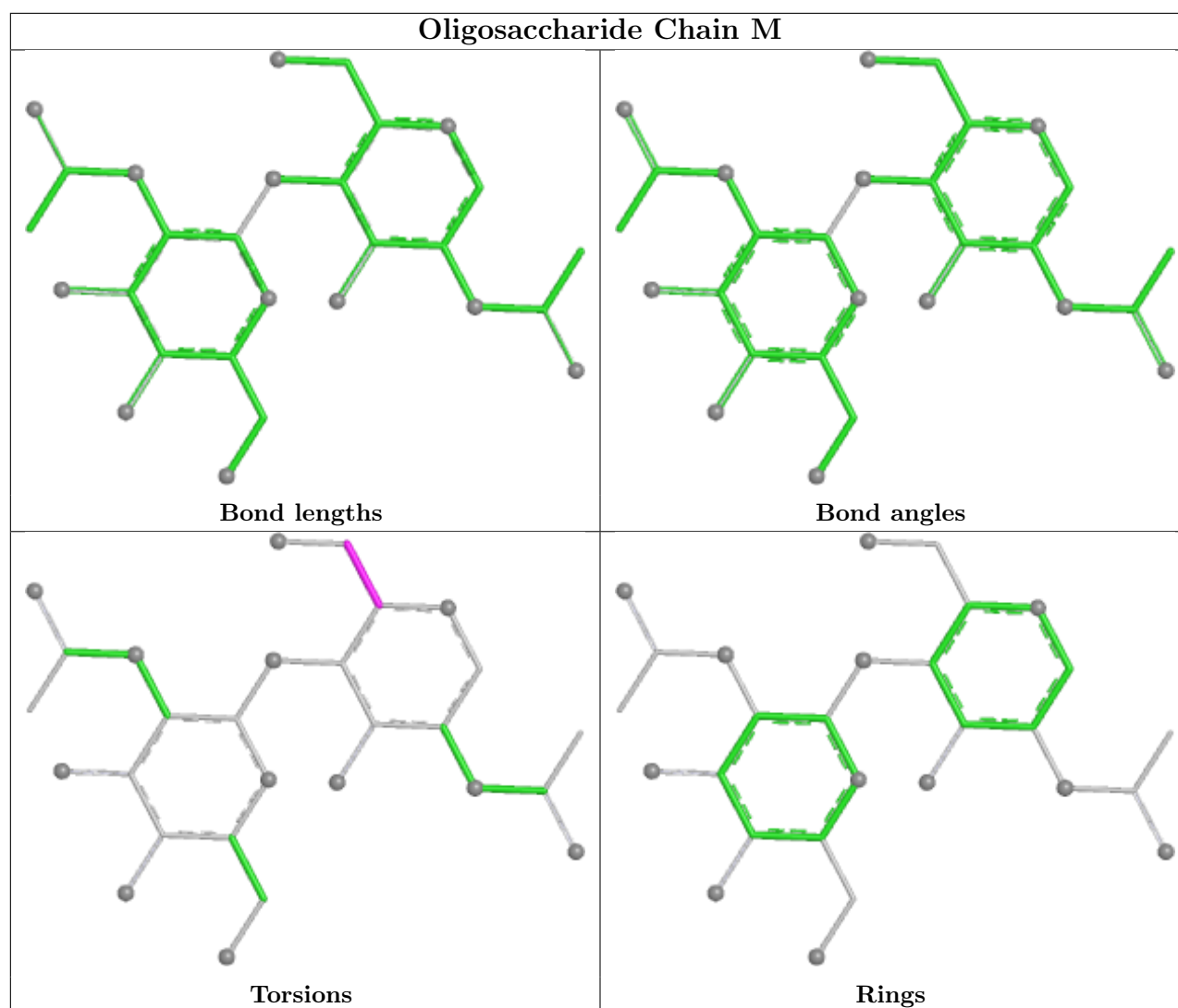
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for oligosaccharide.

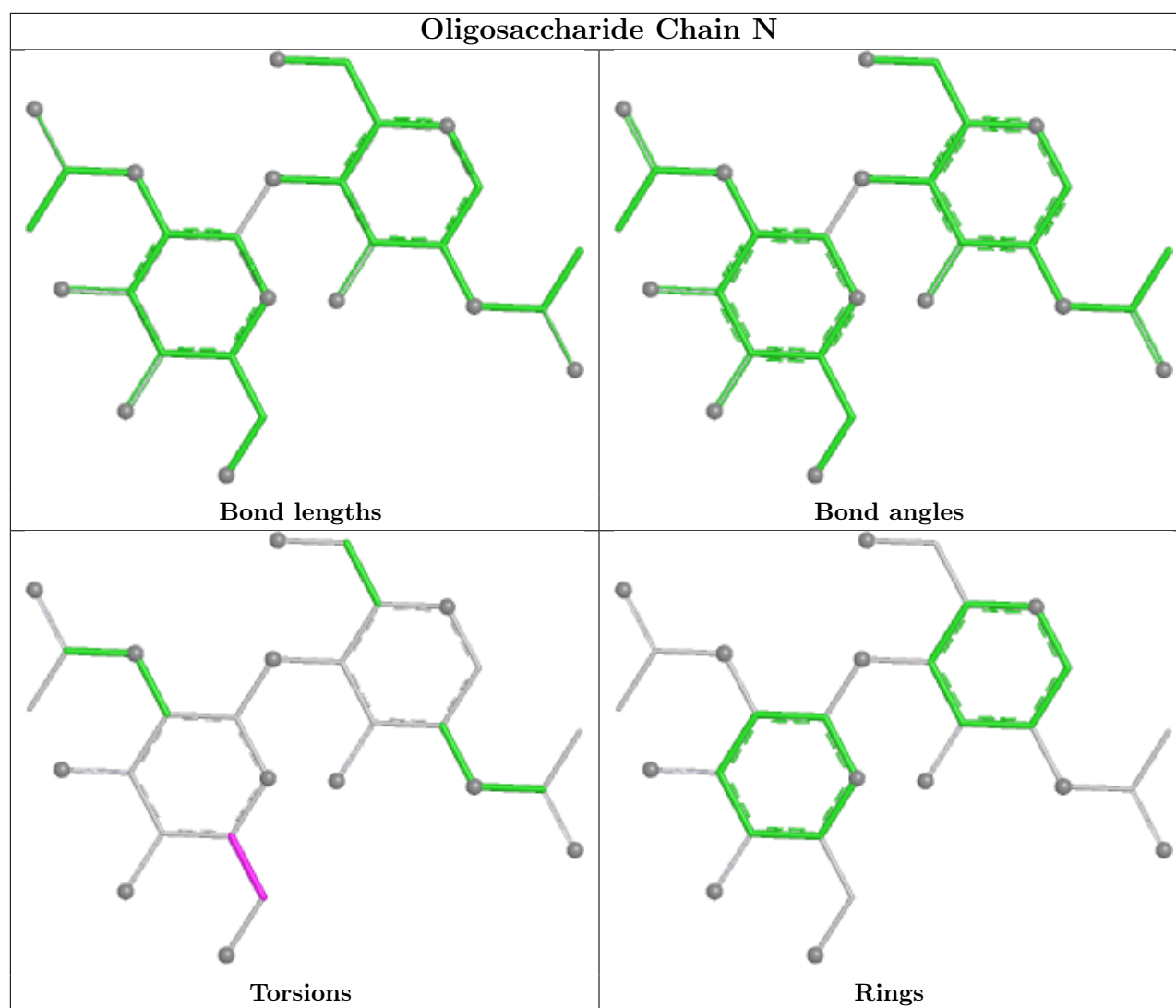


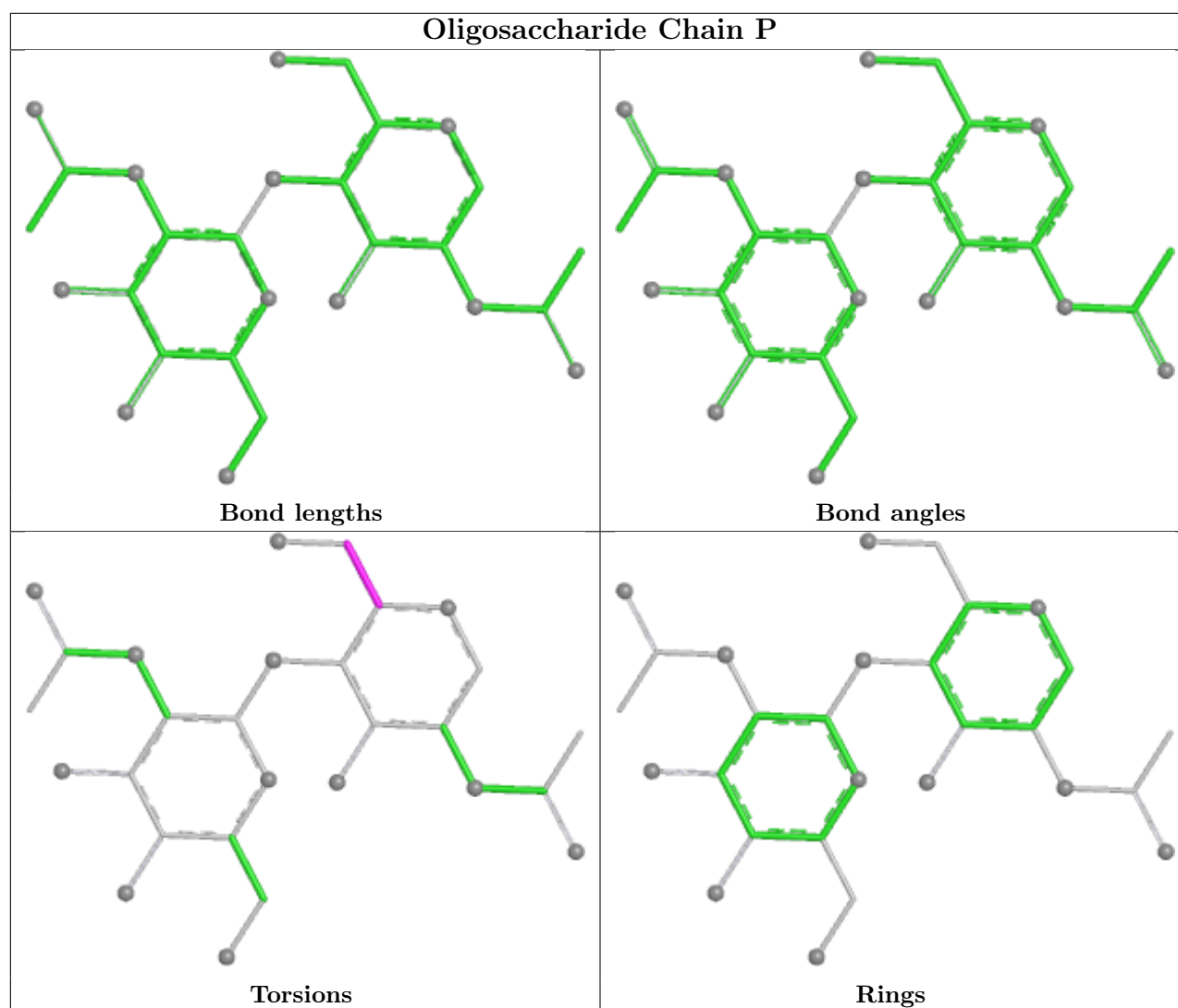


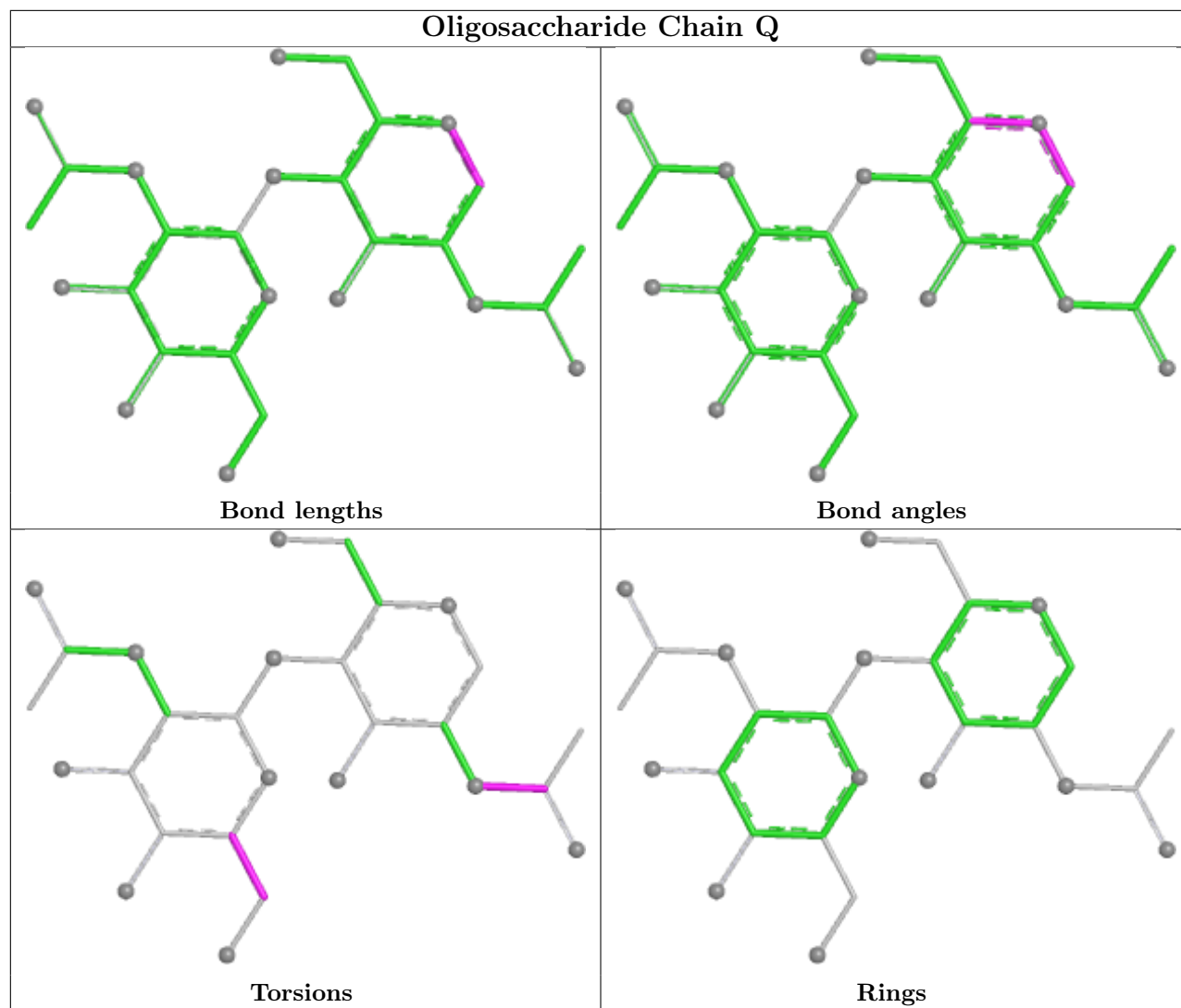


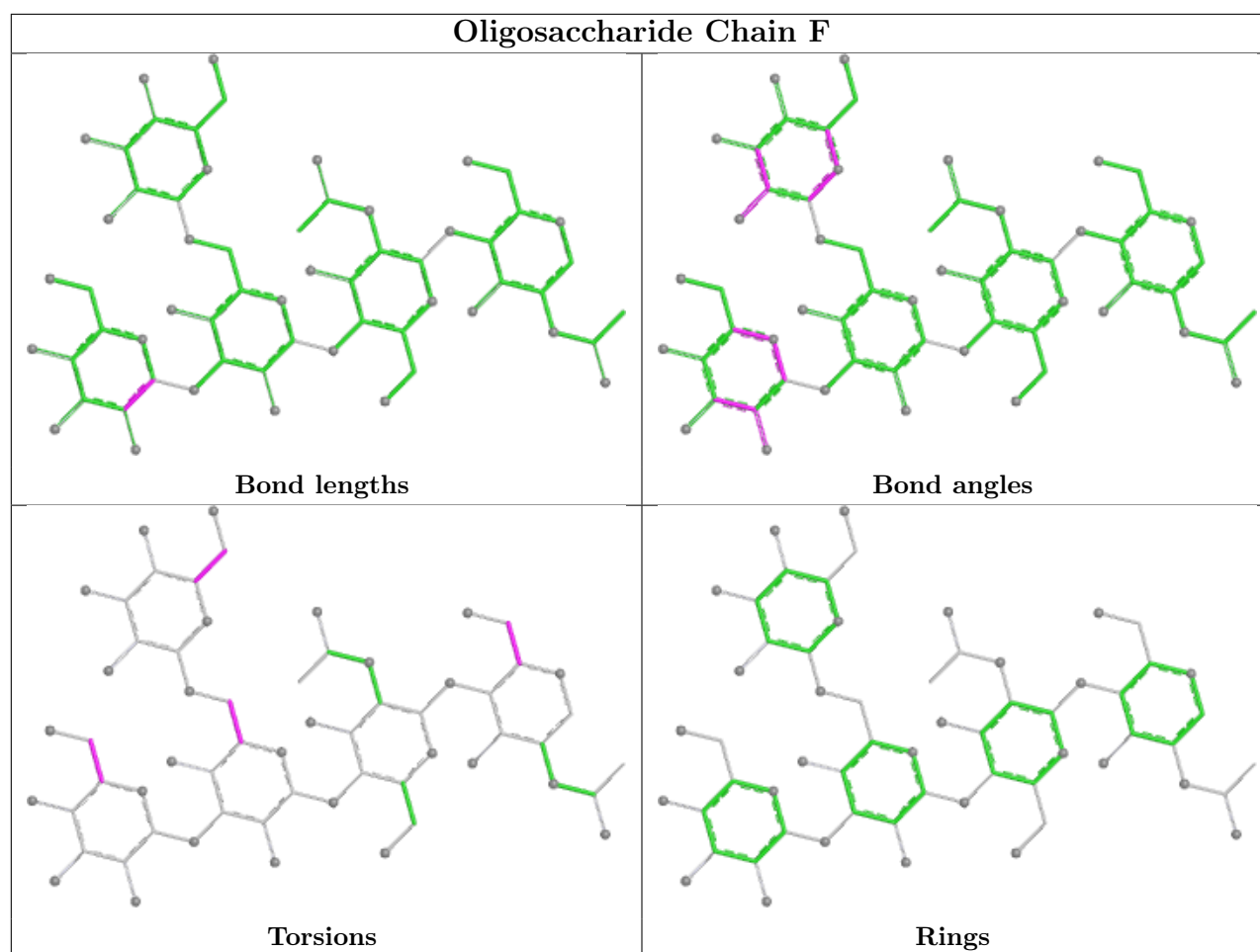


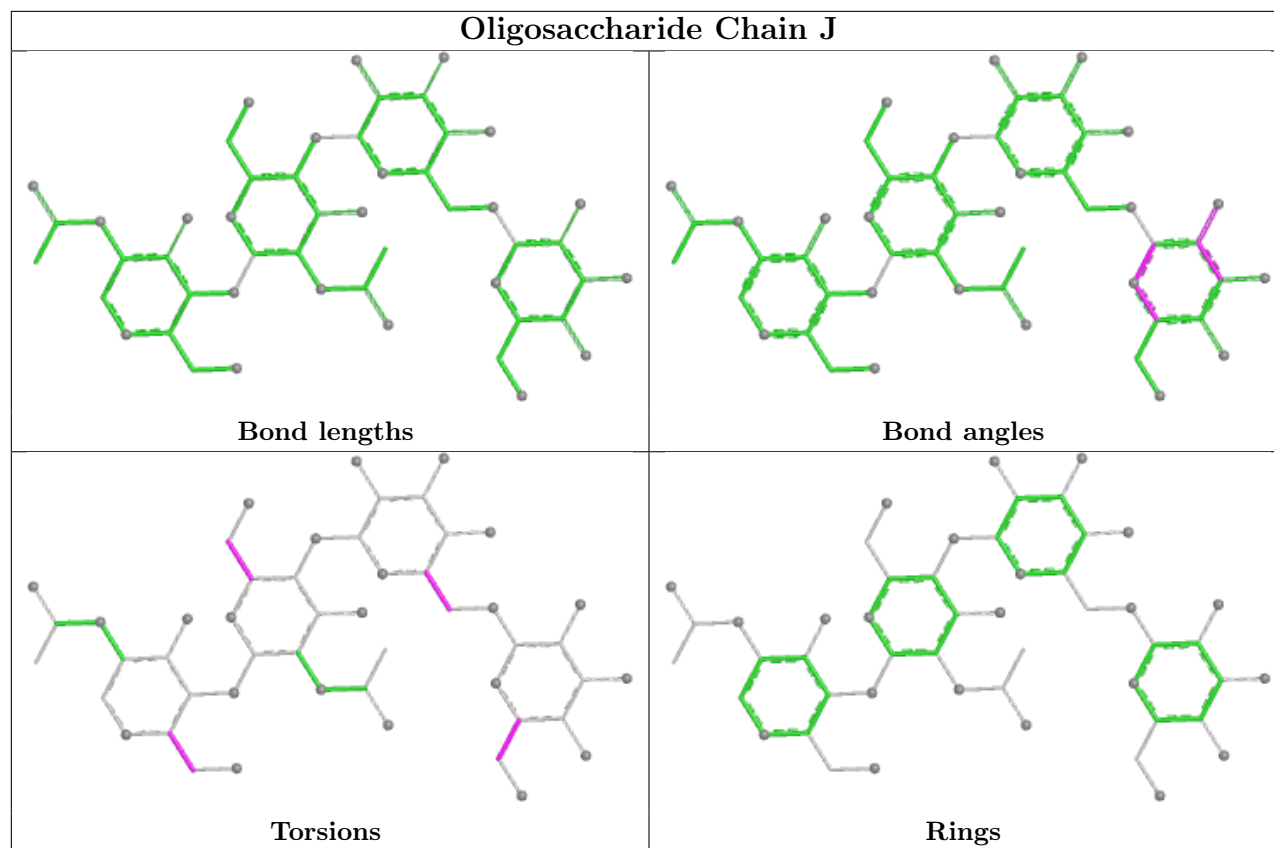


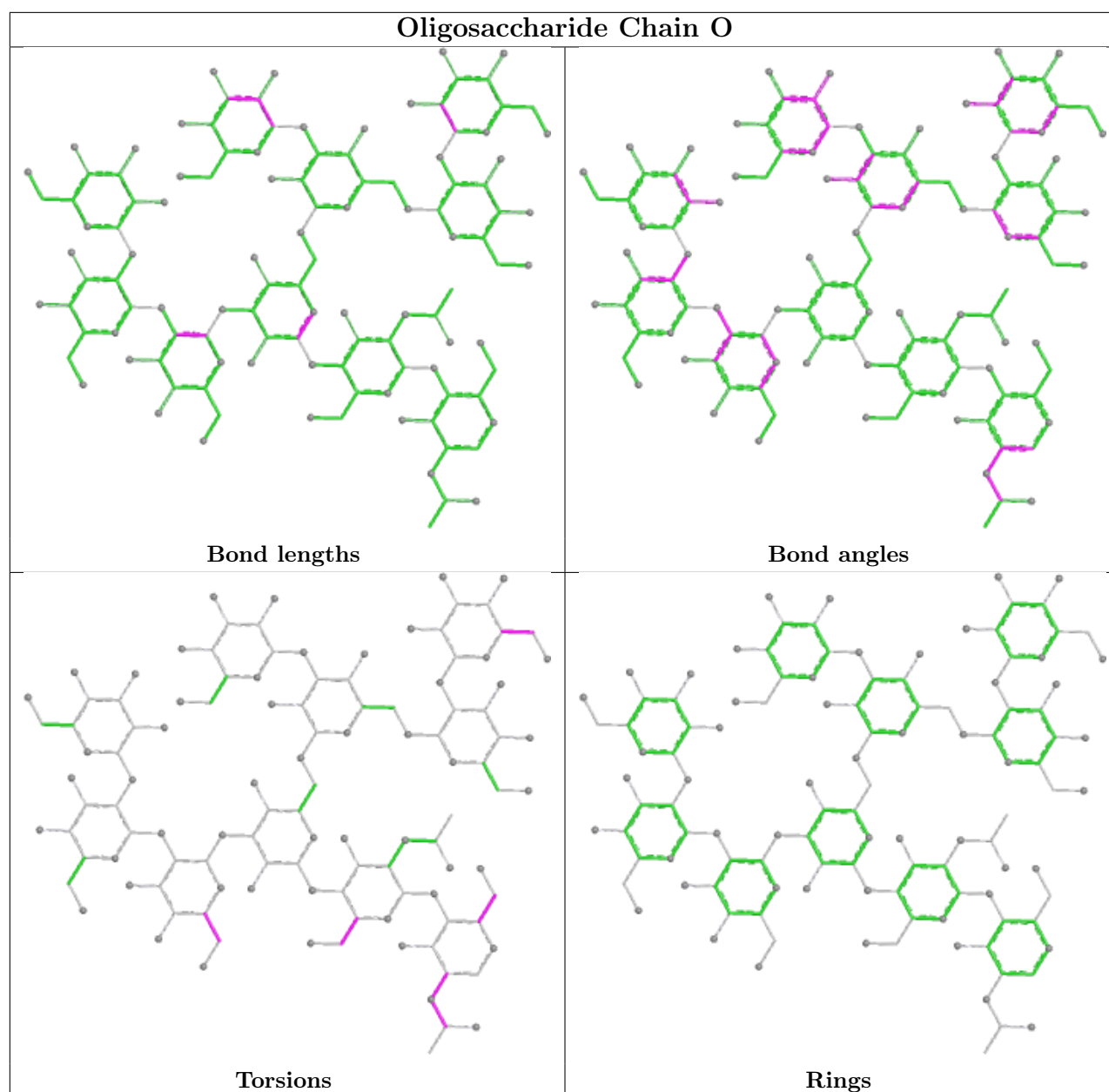












5.6 Ligand geometry ⓘ

7 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
14	NAG	B	701	1	14,14,15	0.37	0	17,19,21	0.59	0
14	NAG	B	702	1	14,14,15	0.30	0	17,19,21	0.45	0
14	NAG	G	614	4	14,14,15	0.29	0	17,19,21	0.49	0
14	NAG	G	638	4	14,14,15	0.21	0	17,19,21	0.45	0
14	NAG	G	617	4	14,14,15	0.24	0	17,19,21	0.47	0
14	NAG	B	703	1	14,14,15	0.29	0	17,19,21	0.47	0
14	NAG	D	201	2	14,14,15	0.52	0	17,19,21	0.51	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
14	NAG	B	701	1	-	1/6/23/26	0/1/1/1
14	NAG	B	702	1	-	2/6/23/26	0/1/1/1
14	NAG	G	614	4	-	2/6/23/26	0/1/1/1
14	NAG	G	638	4	-	3/6/23/26	0/1/1/1
14	NAG	G	617	4	-	2/6/23/26	0/1/1/1
14	NAG	B	703	1	-	1/6/23/26	0/1/1/1
14	NAG	D	201	2	-	3/6/23/26	0/1/1/1

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (14) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
14	G	617	NAG	O5-C5-C6-O6
14	G	614	NAG	O5-C5-C6-O6
14	G	638	NAG	O5-C5-C6-O6
14	G	638	NAG	C4-C5-C6-O6
14	G	617	NAG	C4-C5-C6-O6
14	G	614	NAG	C4-C5-C6-O6
14	B	703	NAG	O5-C5-C6-O6
14	B	702	NAG	C4-C5-C6-O6
14	D	201	NAG	O5-C5-C6-O6
14	B	702	NAG	O5-C5-C6-O6
14	B	701	NAG	O5-C5-C6-O6
14	D	201	NAG	C3-C2-N2-C7

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Mol	Chain	Res	Type	Atoms
14	D	201	NAG	C1-C2-N2-C7
14	G	638	NAG	C1-C2-N2-C7

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
14	D	201	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	B	126/153 (82%)	1.09	29 (23%) 2 2	13, 41, 76, 101	0
2	D	128/153 (83%)	1.94	52 (40%) 0 0	27, 75, 116, 127	0
3	E	110/130 (84%)	1.77	44 (40%) 1 0	31, 59, 103, 120	0
4	G	429/481 (89%)	0.86	58 (13%) 7 6	10, 34, 85, 113	0
5	H	226/244 (92%)	1.17	39 (17%) 4 4	15, 51, 92, 111	0
6	L	210/217 (96%)	0.60	13 (6%) 26 26	16, 39, 66, 91	0
7	U	128/145 (88%)	2.57	84 (65%) 0 0	37, 87, 123, 143	0
8	V	99/122 (81%)	3.44	90 (90%) 0 0	68, 108, 136, 152	0
All	All	1456/1645 (88%)	1.38	409 (28%) 1 1	10, 49, 112, 152	0

All (409) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
8	V	76	SER	7.5
1	B	660	LEU	7.2
4	G	465	THR	7.2
8	V	99	GLY	7.0
8	V	20	THR	6.8
3	E	68	TYR	6.6
7	U	14	PRO	6.5
4	G	505	VAL	6.5
8	V	86	TYR	6.3
2	D	41	ALA	6.3
8	V	48	ILE	6.2
8	V	15	ILE	6.1
3	E	69	TRP	6.0
2	D	109	LEU	6.0
4	G	135	THR	5.9
4	G	188	ASN	5.9

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Mol	Chain	Res	Type	RSRZ
2	D	80	MET	5.9
7	U	7	SER	5.9
4	G	72	HIS	5.8
7	U	8	GLY	5.6
8	V	22	ASN	5.5
6	L	210	THR	5.5
8	V	63	SER	5.4
6	L	184	MET	5.3
4	G	71	THR	5.3
3	E	106(A)	LEU	5.2
8	V	58	VAL	5.2
7	U	91	TYR	5.1
8	V	47	LEU	5.1
8	V	71	PHE	5.0
8	V	80	ALA	4.9
3	E	14	SER	4.9
8	V	104	LEU	4.9
4	G	430	ILE	4.9
7	U	10	ALA	4.9
7	U	36	TRP	4.8
7	U	109	VAL	4.8
1	B	574	LYS	4.7
4	G	151	ARG	4.7
4	G	280	ASN	4.7
2	D	57	LYS	4.6
7	U	90	TYR	4.5
4	G	69	TRP	4.5
5	H	34	TRP	4.5
8	V	54	VAL	4.5
8	V	62	PHE	4.4
2	D	18	VAL	4.4
8	V	70	SER	4.4
7	U	61	GLY	4.4
1	B	548	ILE	4.4
7	U	3	HIS	4.4
2	D	82(C)	LEU	4.4
8	V	31	SER	4.3
8	V	44	PRO	4.3
1	B	663	LEU	4.2
8	V	32	ASP	4.2
3	E	40	PRO	4.2
3	E	41	GLY	4.2

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Mol	Chain	Res	Type	RSRZ
7	U	18	VAL	4.2
8	V	82	ASP	4.2
7	U	15	GLY	4.2
2	D	108	LEU	4.2
8	V	14	SER	4.2
8	V	102	SER	4.2
8	V	73	LEU	4.2
8	V	36	TYR	4.1
7	U	80	MET	4.1
7	U	9	THR	4.1
2	D	88	GLY	4.1
8	V	30	GLY	4.0
8	V	103	ARG	4.0
5	H	78	LEU	4.0
7	U	27	TYR	4.0
7	U	92	CYS	4.0
7	U	40	ALA	3.9
2	D	23	LYS	3.9
4	G	360	ARG	3.9
6	L	49	TYR	3.9
3	E	2	SER	3.9
8	V	37	GLN	3.9
1	B	662	ALA	3.9
8	V	13	VAL	3.9
8	V	3	HIS	3.9
8	V	105	HIS	3.9
8	V	55	GLU	3.8
7	U	45	LEU	3.8
8	V	21	ILE	3.8
8	V	16	GLY	3.8
7	U	5	VAL	3.8
1	B	614	TRP	3.8
4	G	428	GLN	3.8
2	D	17	SER	3.8
2	D	91	PHE	3.8
2	D	5	VAL	3.8
2	D	107	THR	3.8
1	B	648	GLU	3.7
8	V	6	GLN	3.7
8	V	96	GLN	3.7
8	V	81	ASP	3.7
7	U	93	ALA	3.7

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Mol	Chain	Res	Type	RSRZ
7	U	47	TRP	3.7
8	V	75	ILE	3.7
7	U	112	SER	3.7
8	V	67	PHE	3.7
7	U	76(B)	ASP	3.7
3	E	8	ALA	3.7
8	V	53	SER	3.7
5	H	208	LYS	3.7
3	E	107	GLY	3.7
7	U	29	PHE	3.7
4	G	68	VAL	3.7
7	U	16	ALA	3.6
8	V	7	SER	3.6
5	H	51	ILE	3.6
7	U	82	ILE	3.6
2	D	84	SER	3.6
4	G	412	ASP	3.6
4	G	152	GLY	3.6
3	E	11	VAL	3.6
2	D	20	ILE	3.6
2	D	83	THR	3.6
3	E	65	TYR	3.6
8	V	57	GLY	3.5
4	G	211	GLU	3.5
7	U	82(C)	LEU	3.5
5	H	209	VAL	3.5
2	D	44	GLY	3.5
8	V	4	VAL	3.5
8	V	29	VAL	3.4
5	H	56	SER	3.4
7	U	28	THR	3.4
2	D	67	VAL	3.4
1	B	571	TRP	3.4
7	U	82(B)	GLY	3.4
8	V	46	LEU	3.4
8	V	98	PHE	3.4
7	U	103	TRP	3.4
8	V	35	TRP	3.4
3	E	12	SER	3.4
4	G	458	GLY	3.4
7	U	53	GLN	3.4
3	E	15	LEU	3.3

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Mol	Chain	Res	Type	RSRZ
4	G	354	GLY	3.3
7	U	104	GLY	3.3
8	V	24	GLN	3.3
8	V	85	THR	3.3
8	V	19	VAL	3.3
8	V	45	LYS	3.3
8	V	97	PHE	3.3
4	G	499	THR	3.3
8	V	33	LEU	3.3
8	V	89	GLN	3.3
1	B	658	GLN	3.3
6	L	7	TYR	3.3
7	U	39	GLN	3.3
2	D	51	ILE	3.3
2	D	82	ILE	3.3
7	U	73	LEU	3.2
5	H	57	THR	3.2
5	H	19	SER	3.2
7	U	75	GLN	3.2
2	D	37	ILE	3.2
2	D	60	ALA	3.2
5	H	24	VAL	3.2
8	V	72	ASN	3.2
4	G	466	GLU	3.2
3	E	80	PRO	3.2
8	V	27	GLN	3.2
2	D	42	GLY	3.2
7	U	2	ALA	3.2
2	D	29	PHE	3.2
2	D	15	GLY	3.2
8	V	52	SER	3.2
8	V	100	ARG	3.1
2	D	110	THR	3.1
8	V	74	THR	3.1
3	E	99	GLY	3.1
1	B	570	VAL	3.1
7	U	111	VAL	3.1
8	V	87	TYR	3.1
3	E	82	ASP	3.1
8	V	69	THR	3.1
8	V	51	THR	3.1
7	U	26	GLY	3.1

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Mol	Chain	Res	Type	RSRZ
7	U	94	ARG	3.1
3	E	42	ARG	3.1
5	H	187	LEU	3.0
8	V	17	ASP	3.0
7	U	89	VAL	3.0
8	V	78	LEU	3.0
5	H	29	ILE	3.0
7	U	44	GLY	3.0
7	U	107	THR	3.0
4	G	58	ALA	3.0
1	B	615	SER	3.0
4	G	352	HIS	3.0
8	V	34	HIS	3.0
2	D	14	PRO	3.0
8	V	59	PRO	3.0
5	H	158	THR	3.0
3	E	59	SER	3.0
8	V	88	CYS	3.0
8	V	40	PRO	3.0
4	G	335	LYS	3.0
7	U	60	GLY	3.0
7	U	102	ALA	3.0
7	U	110	VAL	2.9
1	B	636	SER	2.9
4	G	457	ASP	2.9
4	G	184	ILE	2.9
7	U	87	THR	2.9
2	D	82(A)	ARG	2.9
8	V	42	ARG	2.9
4	G	67	ASN	2.9
1	B	641	ILE	2.9
5	H	33	TYR	2.9
5	H	189	THR	2.9
2	D	16	SER	2.9
5	H	25	SER	2.9
7	U	17	SER	2.9
2	D	68	ASN	2.9
1	B	659	ASP	2.9
8	V	11	LEU	2.9
8	V	90	VAL	2.9
7	U	76(F)	TRP	2.8
4	G	425	ASN	2.8

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Mol	Chain	Res	Type	RSRZ
8	V	49	HIS	2.8
5	H	183	PRO	2.8
2	D	43	ARG	2.8
7	U	105	GLN	2.8
8	V	12	SER	2.8
3	E	102	THR	2.8
7	U	88	ALA	2.8
4	G	456	ARG	2.8
5	H	186	SER	2.8
8	V	77	ASP	2.8
3	E	87	TYR	2.8
2	D	92	CYS	2.8
7	U	76(D)	PRO	2.8
1	B	573	ILE	2.8
5	H	69	ILE	2.8
8	V	91	LEU	2.8
4	G	70	ALA	2.8
3	E	62	PHE	2.8
2	D	38	ARG	2.8
2	D	45	PRO	2.7
8	V	10	SER	2.7
7	U	76(E)	ASP	2.7
7	U	108	THR	2.7
8	V	38	HIS	2.7
3	E	79	ARG	2.7
6	L	53	ASP	2.7
3	E	106	VAL	2.7
2	D	56	ASP	2.7
4	G	387	THR	2.7
7	U	37	PHE	2.7
2	D	8	GLY	2.7
1	B	613	SER	2.7
7	U	100(B)	SER	2.7
3	E	43	ALA	2.7
3	E	85	THR	2.7
7	U	100(D)	ALA	2.7
2	D	49	GLY	2.6
7	U	57	VAL	2.6
2	D	12	THR	2.6
2	D	70	THR	2.6
8	V	79	GLN	2.6
5	H	136	LEU	2.6

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Mol	Chain	Res	Type	RSRZ
7	U	38	ARG	2.6
7	U	83	LYS	2.6
2	D	21	SER	2.6
6	L	18	THR	2.6
4	G	359	ILE	2.6
3	E	3	VAL	2.6
8	V	43	ALA	2.6
4	G	79	PRO	2.6
8	V	9	SER	2.5
7	U	76(C)	ASP	2.5
4	G	284	ILE	2.5
5	H	132	GLY	2.5
7	U	20	VAL	2.5
2	D	10	THR	2.5
3	E	27	ASN	2.5
7	U	12	LYS	2.5
1	B	638	TYR	2.5
5	H	55	GLU	2.5
5	H	77	GLN	2.5
3	E	84	THR	2.5
5	H	110	SER	2.5
3	E	78	LEU	2.5
8	V	18	ARG	2.5
7	U	46	GLU	2.5
4	G	432	GLN	2.5
5	H	71	VAL	2.5
5	H	112	ALA	2.5
2	D	19	LYS	2.5
8	V	84	ALA	2.5
3	E	70	SER	2.4
6	L	60	GLU	2.4
7	U	79	TYR	2.4
4	G	281	ALA	2.4
7	U	35	PHE	2.4
7	U	82(A)	ARG	2.4
4	G	349	LEU	2.4
6	L	67(A)	ILE	2.4
7	U	99	GLY	2.4
8	V	8	PRO	2.4
6	L	185	GLN	2.4
4	G	397	SER	2.4
7	U	74	SER	2.4

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Mol	Chain	Res	Type	RSRZ
7	U	100	ASP	2.4
7	U	76(G)	GLY	2.4
2	D	62	ALA	2.4
4	G	467	THR	2.4
2	D	81	GLU	2.4
3	E	105	SER	2.4
4	G	153	GLU	2.4
2	D	47	TRP	2.4
3	E	103	LYS	2.4
7	U	48	VAL	2.4
5	H	68	ILE	2.3
8	V	5	THR	2.3
3	E	29	SER	2.3
5	H	28	SER	2.3
5	H	52	SER	2.3
7	U	97	SER	2.3
4	G	35	TRP	2.3
7	U	96	ARG	2.3
7	U	63	PHE	2.3
3	E	30	HIS	2.3
4	G	357	THR	2.3
1	B	656	ASN	2.3
7	U	86	ASP	2.3
4	G	282	LYS	2.3
6	L	70	ARG	2.3
7	U	41	PRO	2.3
3	E	16	GLY	2.3
6	L	68	GLY	2.3
7	U	6	GLN	2.3
7	U	49	GLY	2.3
7	U	100(E)	LEU	2.3
3	E	48	ILE	2.3
8	V	50	HIS	2.3
8	V	68	HIS	2.3
6	L	188	MET	2.3
1	B	547	GLY	2.3
7	U	50	TRP	2.3
1	B	518	VAL	2.3
5	H	100(R)	VAL	2.3
5	H	174	TYR	2.2
4	G	429	ARG	2.2
1	B	649	SER	2.2

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Mol	Chain	Res	Type	RSRZ
5	H	180	VAL	2.2
5	H	138	CYS	2.2
1	B	657	GLU	2.2
4	G	198	THR	2.2
5	H	31	ASN	2.2
7	U	43	ARG	2.2
8	V	41	GLY	2.2
1	B	655	LYS	2.2
4	G	73	ALA	2.2
1	B	639	THR	2.2
4	G	504	ARG	2.2
6	L	9	ARG	2.2
1	B	645	LEU	2.2
2	D	26	GLY	2.2
7	U	77	ILE	2.2
5	H	156	ALA	2.1
2	D	46	GLU	2.1
7	U	1	ARG	2.1
2	D	22	CYS	2.1
4	G	81	PRO	2.1
3	E	17	GLN	2.1
4	G	183	GLN	2.1
4	G	332	ASN	2.1
5	H	157	LEU	2.1
1	B	546	SER	2.1
3	E	27(B)	VAL	2.1
4	G	333	VAL	2.1
3	E	81	GLU	2.1
1	B	609	PRO	2.1
3	E	88	CYS	2.1
2	D	48	MET	2.1
2	D	90	TYR	2.1
4	G	390	LEU	2.1
5	H	100(O)	TYR	2.1
4	G	431	GLY	2.1
8	V	28	GLY	2.1
8	V	101	GLY	2.1
1	B	651	ASN	2.1
5	H	197	ASN	2.1
3	E	47	ILE	2.1
3	E	104	VAL	2.1
4	G	210	PHE	2.1

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Mol	Chain	Res	Type	RSRZ
8	V	56	ASP	2.1
5	H	36	TRP	2.1
7	U	11	MET	2.1
2	D	39	GLN	2.1
3	E	98	PHE	2.1
3	E	7	SER	2.0
4	G	56	SER	2.0
4	G	74	CYS	2.0
2	D	63	PHE	2.0
8	V	39	LYS	2.0
4	G	470	PRO	2.0
5	H	181	THR	2.0
4	G	111	LEU	2.0
1	B	616	ASN	2.0
4	G	235	GLY	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
9	NAG	A	1	14/15	-	-	17,24,30,35	0
9	NAG	A	2	14/15	-	-	24,33,51,63	0
9	BMA	A	3	11/12	-	-	38,43,51,85	0
9	MAN	A	4	11/12	-	-	79,88,107,114	0
9	MAN	A	5	11/12	-	-	85,95,100,105	0
9	MAN	A	6	11/12	-	-	31,42,50,58	0
10	NAG	C	1	14/15	-	-	87,108,130,143	0
10	NAG	C	2	14/15	-	-	114,146,158,165	0
12	NAG	J	2	14/15	0.31	0.28	41,61,102,104	0
10	NAG	N	2	14/15	0.51	0.26	77,87,116,135	0
10	NAG	P	2	14/15	0.64	0.22	48,88,101,106	0
13	NAG	O	2	14/15	0.68	0.16	32,57,74,82	0
10	NAG	N	1	14/15	0.73	0.19	28,43,64,65	0

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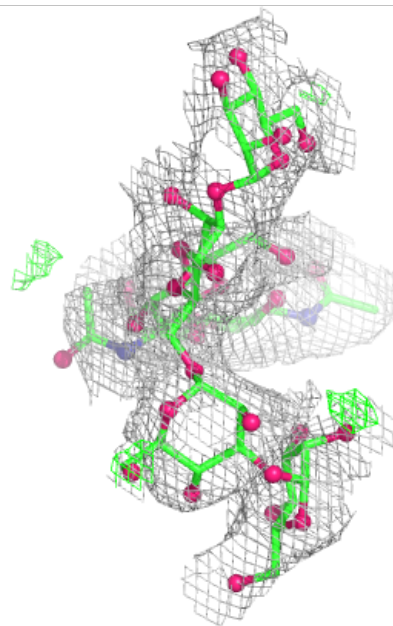
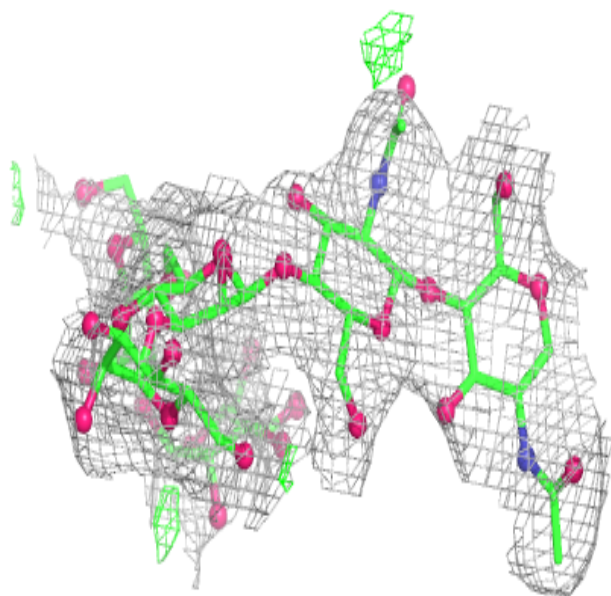
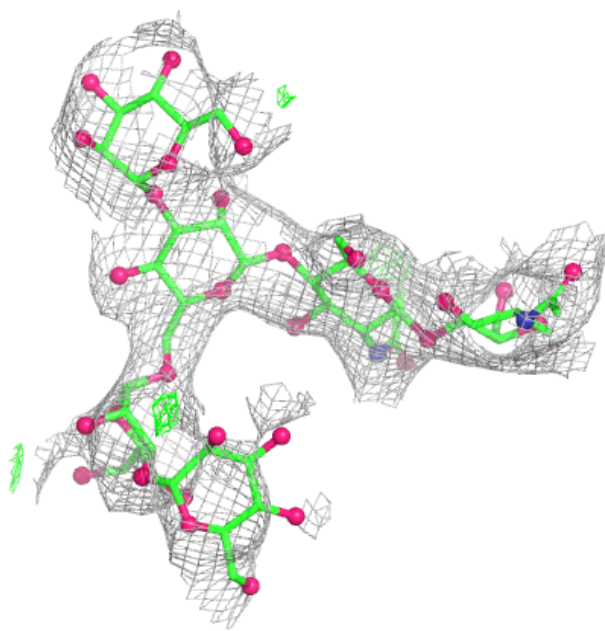
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
12	NAG	J	1	14/15	0.74	0.17	12,26,49,51	0
10	NAG	Q	1	14/15	0.76	0.20	54,68,94,108	0
10	NAG	K	2	14/15	0.79	0.15	100,137,149,149	0
10	NAG	M	2	14/15	0.83	0.13	77,92,117,120	0
10	NAG	Q	2	14/15	0.84	0.15	84,114,126,146	0
10	NAG	K	1	14/15	0.89	0.13	67,98,117,128	0
13	NAG	O	1	14/15	0.92	0.09	33,58,71,77	0
11	NAG	F	1	14/15	-	-	24,56,69,73	0
11	NAG	F	2	14/15	-	-	63,80,113,120	0
11	BMA	F	3	11/12	-	-	123,127,141,142	0
11	MAN	F	4	11/12	-	-	94,119,128,131	0
11	MAN	F	5	11/12	-	-	54,111,124,128	0
10	NAG	I	1	14/15	0.92	0.10	75,96,103,114	0
10	NAG	P	1	14/15	0.93	0.12	46,58,70,94	0
12	BMA	J	3	11/12	-	-	86,107,130,134	0
12	MAN	J	4	11/12	-	-	115,128,149,153	0
10	NAG	M	1	14/15	0.94	0.09	36,52,76,81	0
10	NAG	I	2	14/15	0.94	0.11	72,103,117,121	0
13	BMA	O	3	11/12	-	-	20,39,45,50	0
13	MAN	O	4	11/12	-	-	25,29,39,44	0
13	MAN	O	5	11/12	-	-	22,32,41,44	0
13	MAN	O	6	11/12	-	-	30,37,43,47	0
13	MAN	O	7	11/12	-	-	37,49,66,80	0
13	MAN	O	8	11/12	-	-	42,59,71,72	0
13	MAN	O	9	11/12	-	-	79,86,103,103	0
13	MAN	O	10	11/12	-	-	37,51,77,85	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.

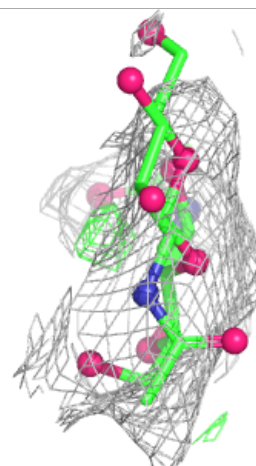
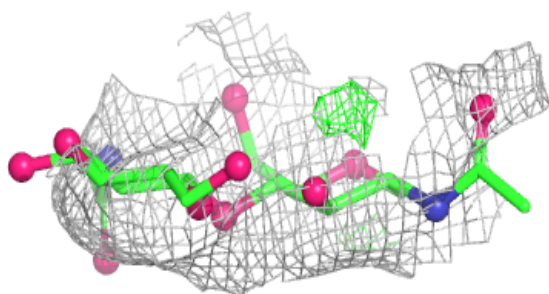
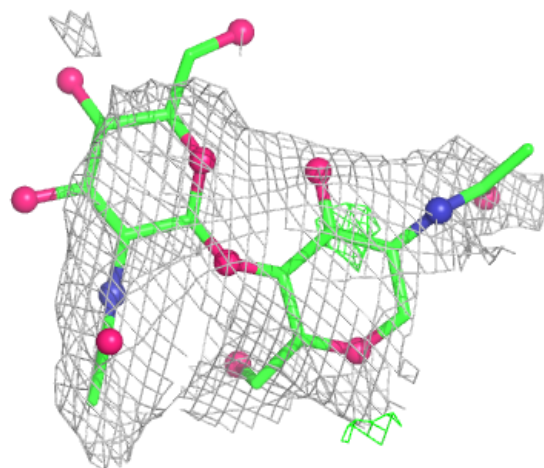
Electron density around Chain A:

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



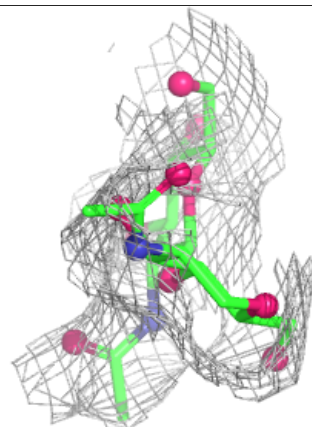
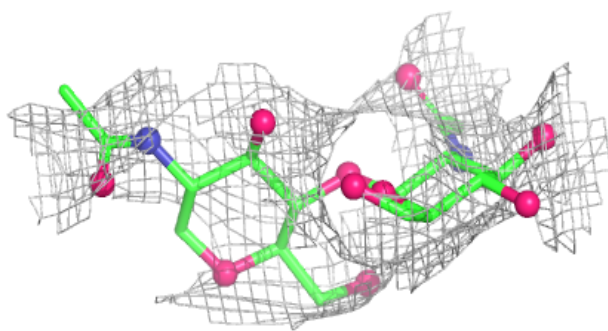
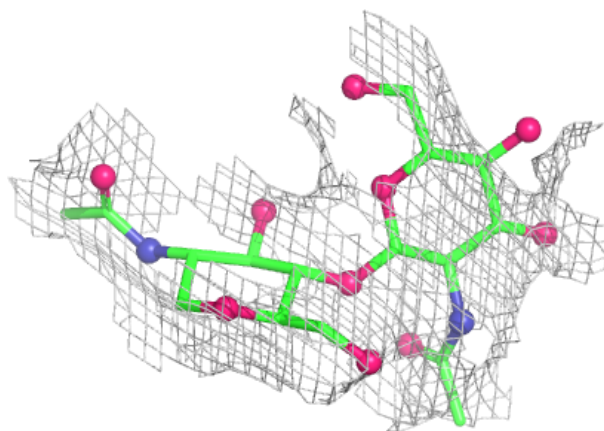
Electron density around Chain C:

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



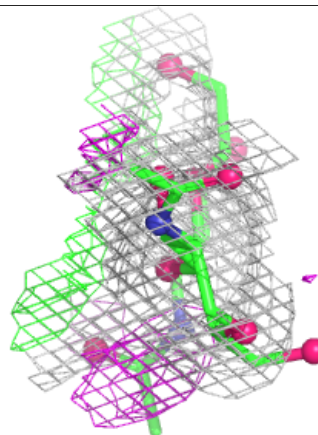
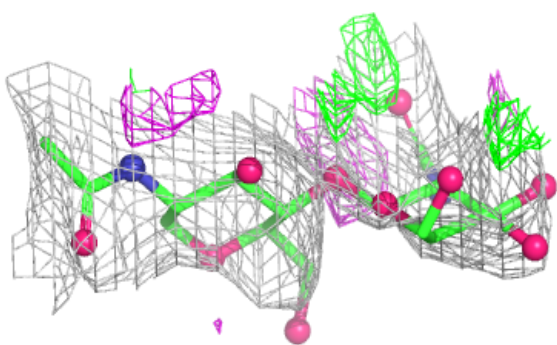
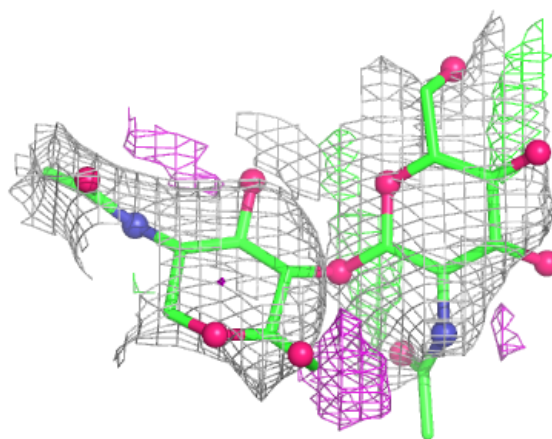
Electron density around Chain I:

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



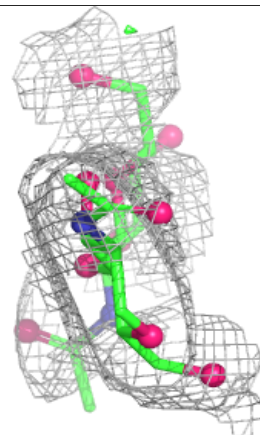
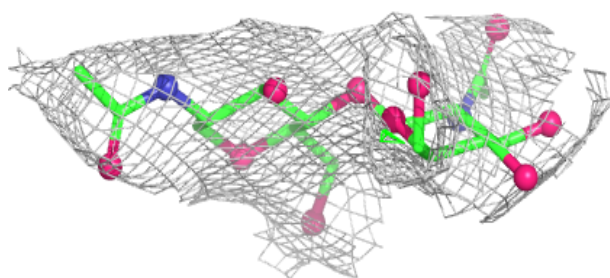
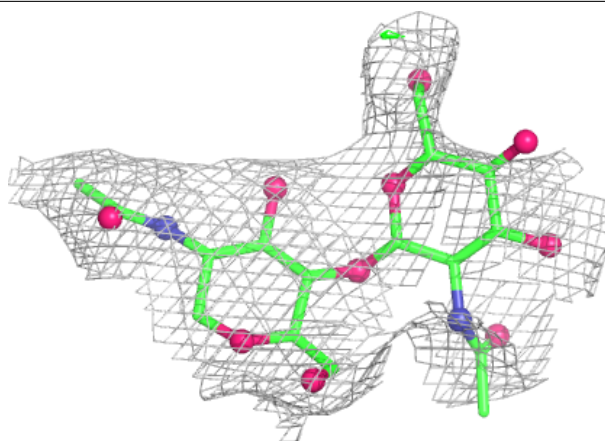
Electron density around Chain K:

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



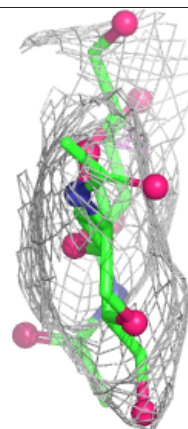
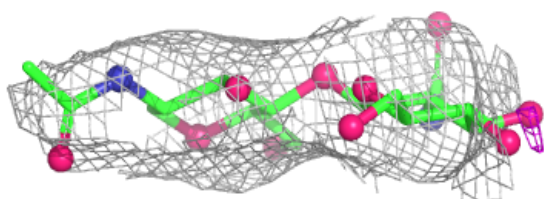
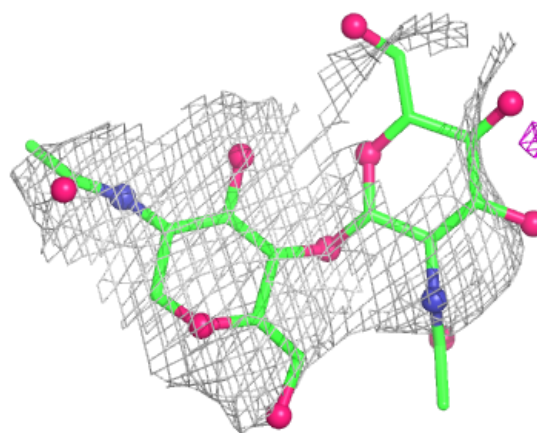
Electron density around Chain M:

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



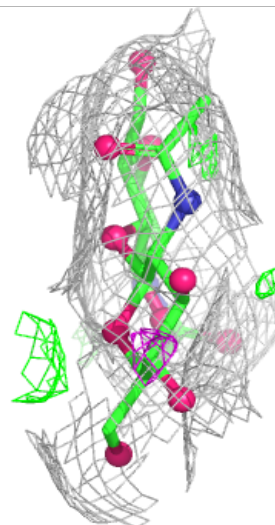
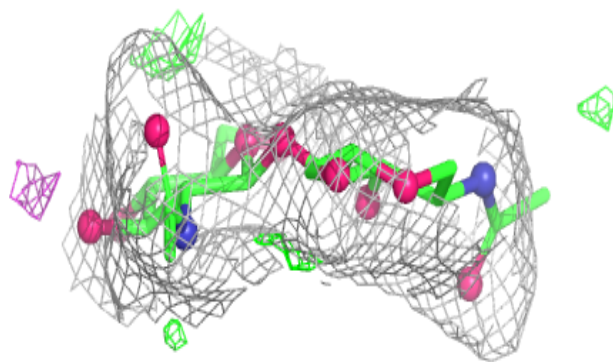
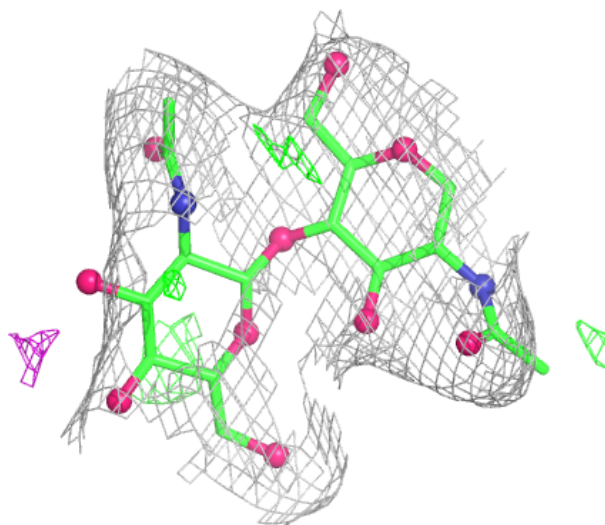
Electron density around Chain N:

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



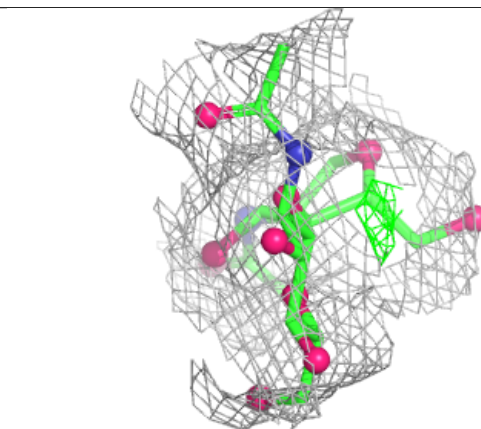
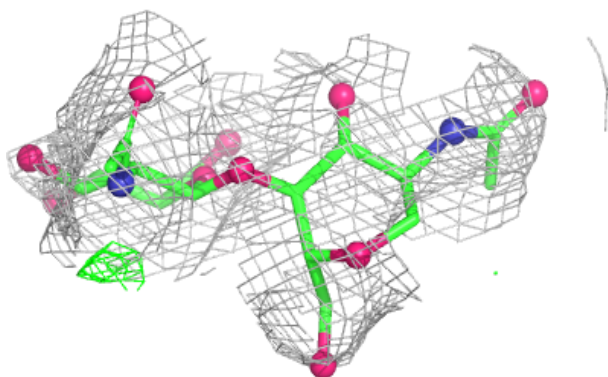
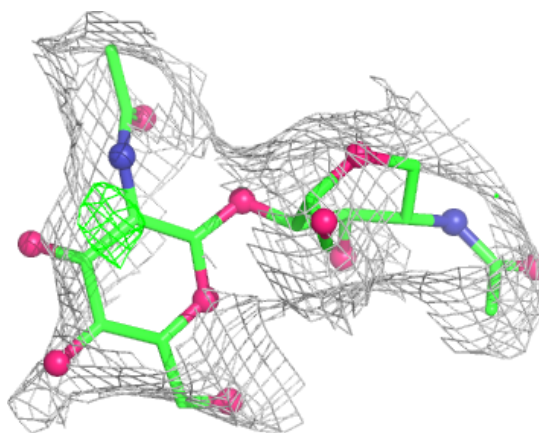
Electron density around Chain P:

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

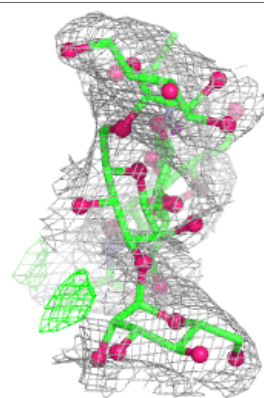
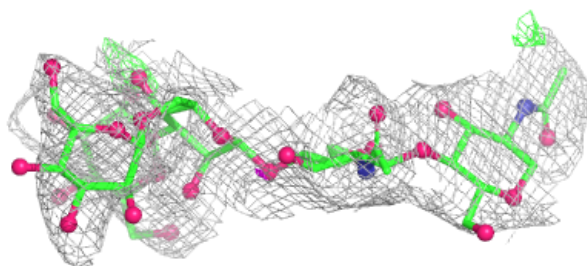
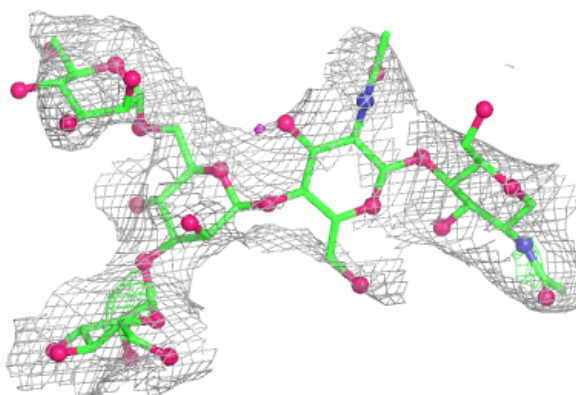


Electron density around Chain Q:

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

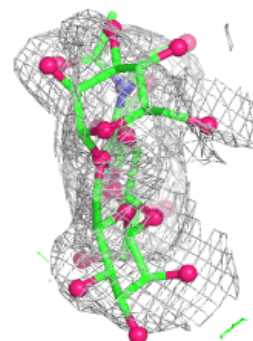
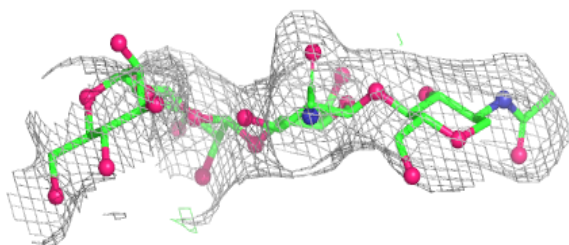
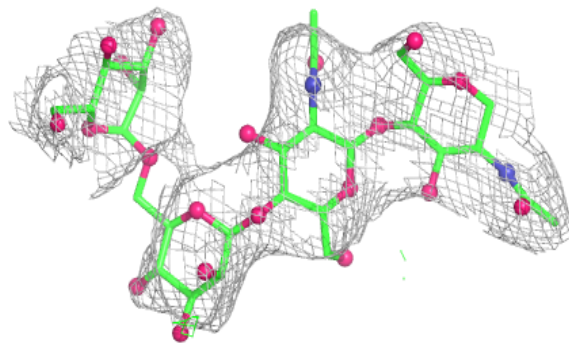
**Electron density around Chain F:**

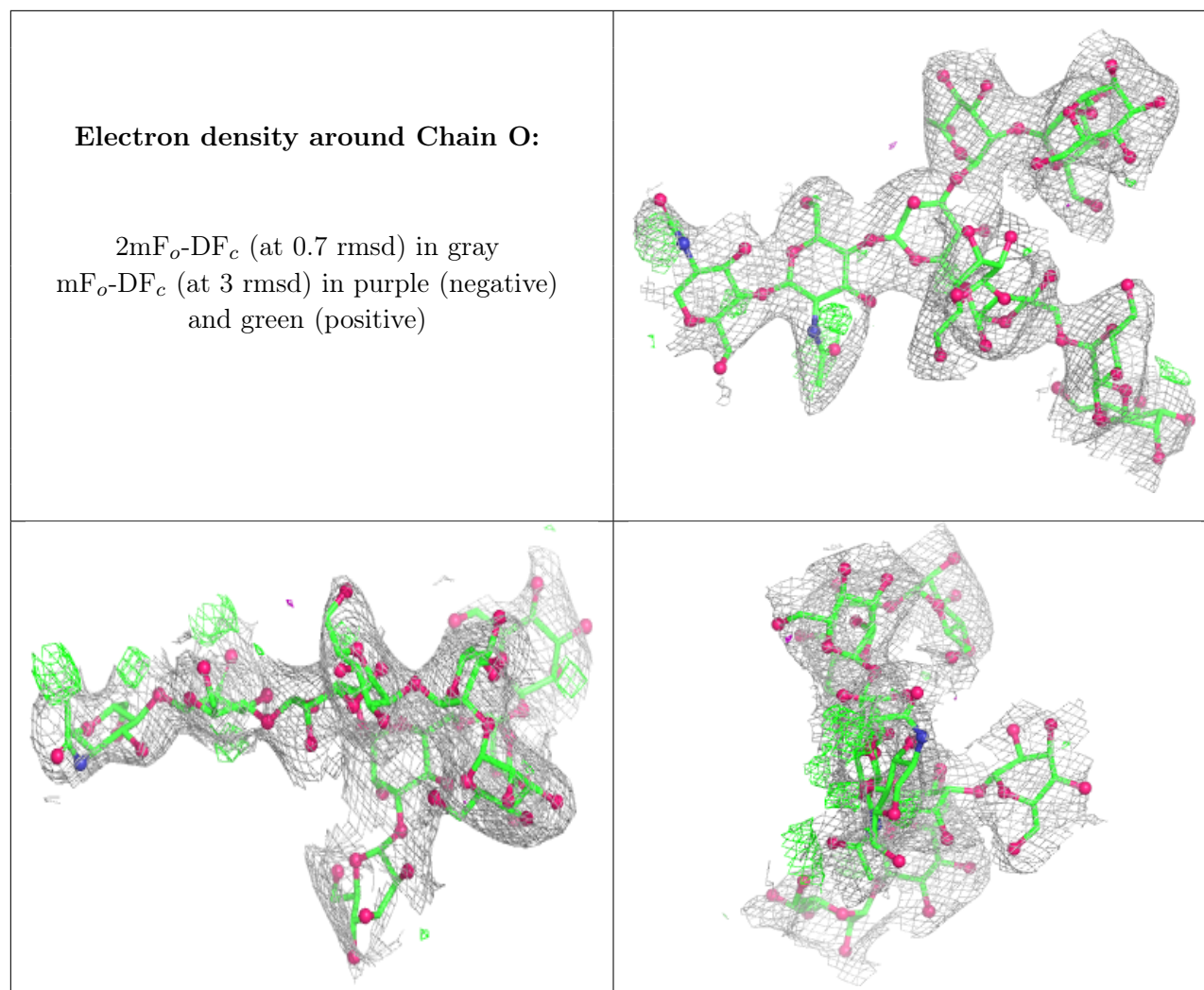
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around Chain J:

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





6.4 Ligands ⓘ

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
14	NAG	B	702	14/15	0.55	0.21	58,96,101,102	0
14	NAG	B	701	14/15	0.63	0.21	85,111,128,139	0
14	NAG	G	617	14/15	0.69	0.23	68,91,99,99	0
14	NAG	B	703	14/15	0.77	0.18	67,103,116,120	0
14	NAG	D	201	14/15	0.81	0.16	64,106,122,128	0
14	NAG	G	614	14/15	0.88	0.11	47,70,96,99	0
14	NAG	G	638	14/15	0.89	0.12	48,57,67,74	0

6.5 Other polymers [i](#)

There are no such residues in this entry.