Full wwPDB X-ray Structure Validation Report

Aug 9, 2020 – 05:23 AM BST

PDB ID	:	6NNJ
Title	:	Crystal Structure of HIV-1 BG505 SOSIP.664 Prefusion Env Trimer Bound to CH31 scFv in Complex with Crystallization Chaperones 3H109L Fab and 35O22 scFv at 3.1 Angstrom
Authors	:	Lai, Y.-T.; Kwong, P.D.
Deposited on	:	2019-01-15
Resolution	:	2.60 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 (i) symbol.

The following versions of software and data (see references (1)) were used in the production of this report:

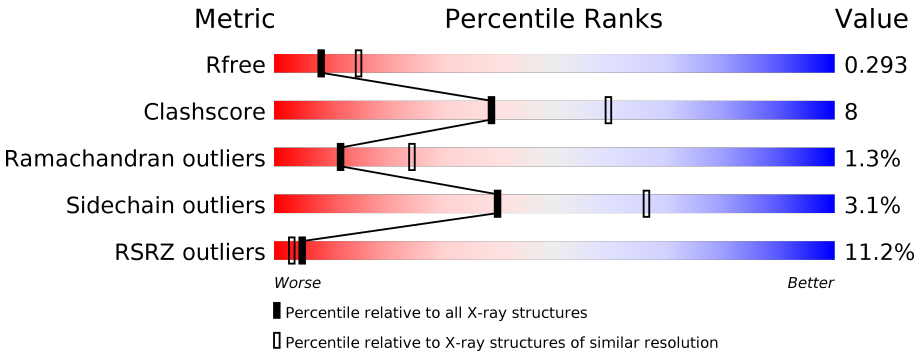
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.13.1
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry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

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.60 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	3163 (2.60-2.60)
Clashscore	141614	3518 (2.60-2.60)
Ramachandran outliers	138981	3455 (2.60-2.60)
Sidechain outliers	138945	3455 (2.60-2.60)
RSRZ outliers	127900	3104 (2.60-2.60)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for >=3, 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions <=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	69% 14% • 17%
2	D	153	24% 60% 22% • 16%
3	E	130	9% 69% 15% 15%
4	G	481	% 67% 21% • 11%
5	H	244	9% 73% 18% • 7%
6	L	217	2% 74% 21% • •

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Mol	Chain	Length	Quality of chain
7	U	137	29% 85% 8% 6%
8	V	117	35% 76% 10% 14%
9	A	6	33% 17% 50%
10	C	2	100%
10	J	2	50% 50%
10	K	2	100%
10	M	2	100%
10	O	2	50% 50%
11	F	5	60% 40%
12	I	4	50% 50%
13	N	10	10% 70% 20%

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
10	NAG	C	2	-	-	-	X

2 Entry composition

There are 14 unique types of molecules in this entry. The entry contains 11874 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	127	Total	C	N	O	S	0	0	0
			1009	641	173	189	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	430	Total	C	N	O	S	0	0	0
			3397	2140	602	628	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			

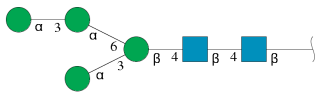
- Molecule 7 is a protein called CH31 scFv heavy chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
7	U	129	Total	C	N	O	S	0	0	0
			1027	652	184	187	4			

- Molecule 8 is a protein called CH31 scFv light chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
8	V	101	Total	C	N	O	S	0	0	0
			761	472	130	156	3			

- 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.

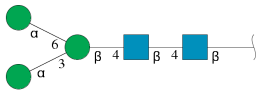


Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
9	A	6	Total	C	N	O	0	0	0
			72	40	2	30			

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

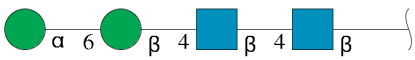
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
10	C	2	Total	C	N	O	0	0	0
			28	16	2	10			
10	J	2	Total	C	N	O	0	0	0
			28	16	2	10			
10	K	2	Total	C	N	O	0	0	0
			28	16	2	10			
10	M	2	Total	C	N	O	0	0	0
			28	16	2	10			
10	O	2	Total	C	N	O	0	0	0
			28	16	2	10			

- 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.



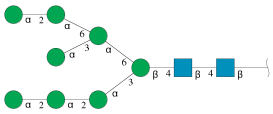
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
11	F	5	Total	C	N	O	0	0	0
			61	34	2	25			

- 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.



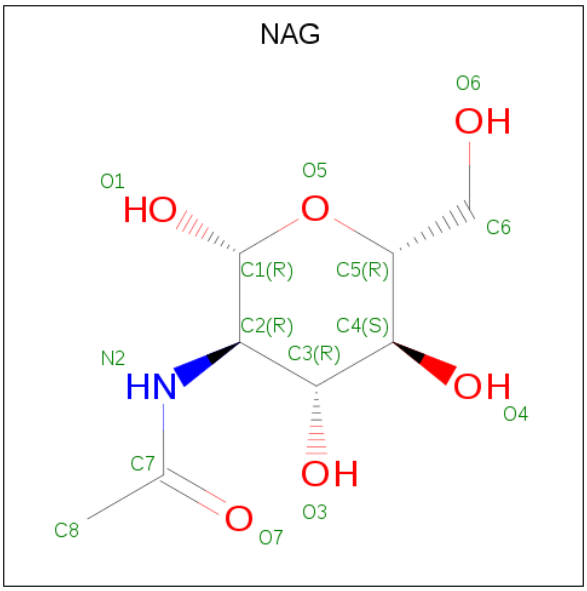
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf	Trace
12	I	4	Total C N O 50 28 2 20	0	0	0

- 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	N	10	Total C N O 116 64 2 50	0	0	0

- Molecule 14 is 2-acetamido-2-deoxy-beta-D-glucopyranose (three-letter code: NAG) (formula: $C_8H_{15}NO_6$).



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

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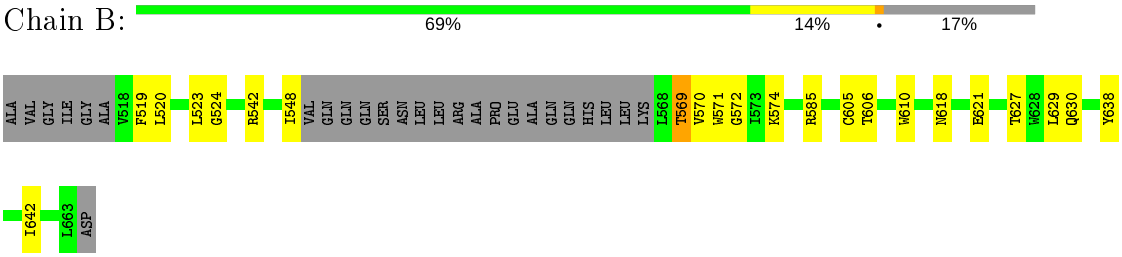
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
14	G	1	Total	C	N	O	0	0
			14	8	1	5		
14	G	1	Total	C	N	O	0	0
			14	8	1	5		
14	G	1	Total	C	N	O	0	0
			14	8	1	5		
14	G	1	Total	C	N	O	0	0
			14	8	1	5		

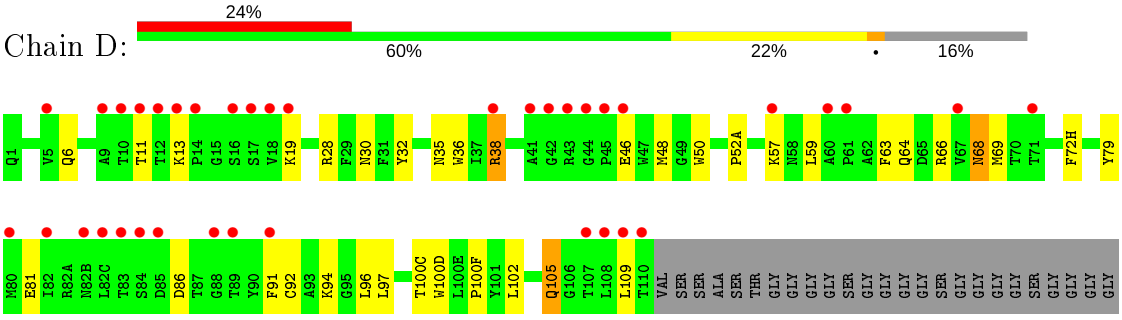
3 Residue-property plots (i)

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.

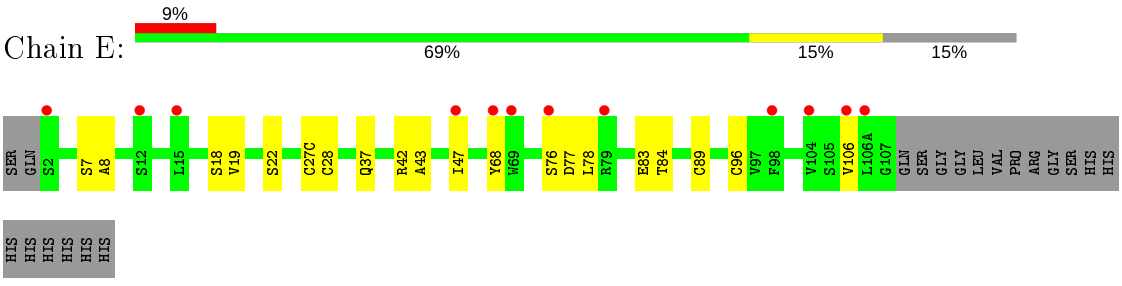
- Molecule 1: Envelope glycoprotein gp41



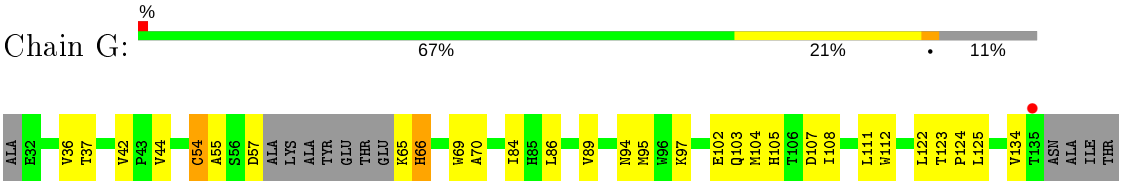
- Molecule 2: 35O22 scFv heavy chain

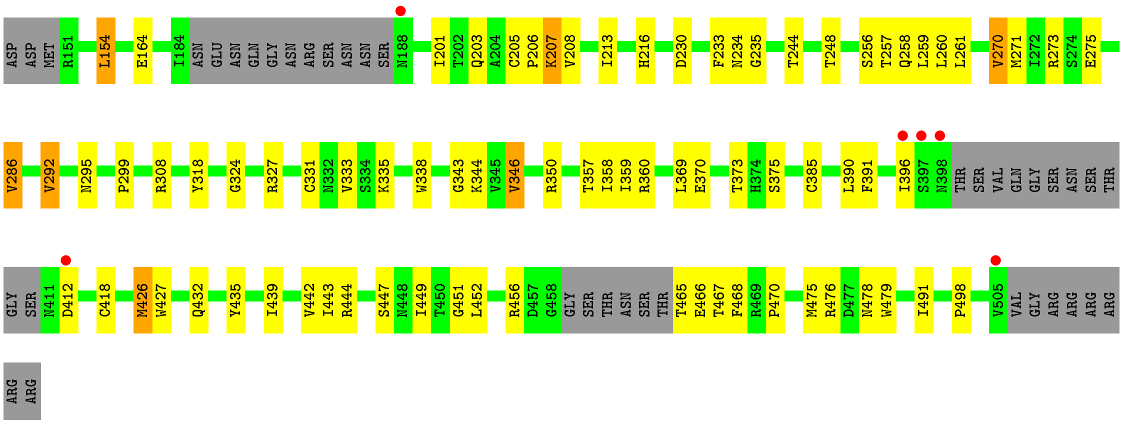


- Molecule 3: 35O22 scFv light chain

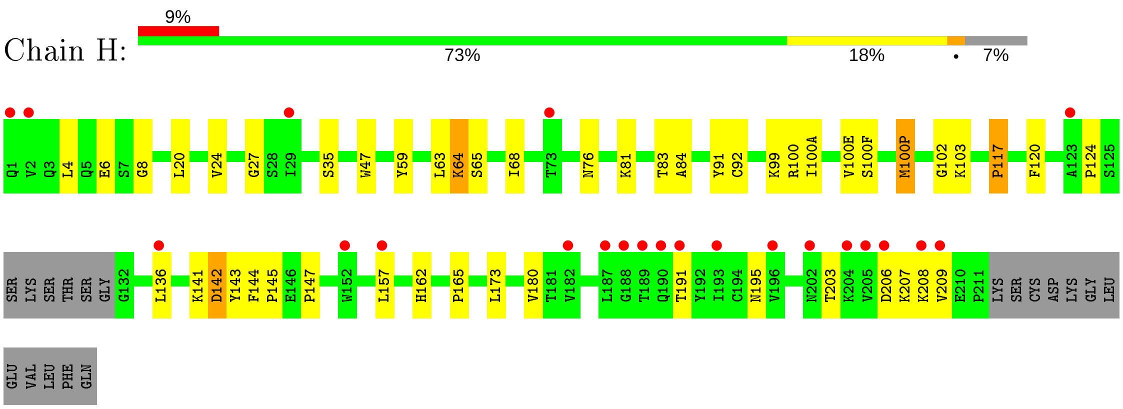


- Molecule 4: Envelope glycoprotein gp120

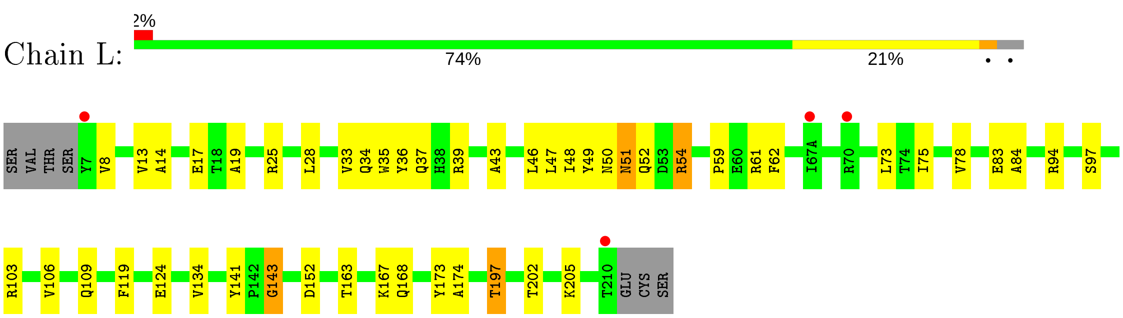




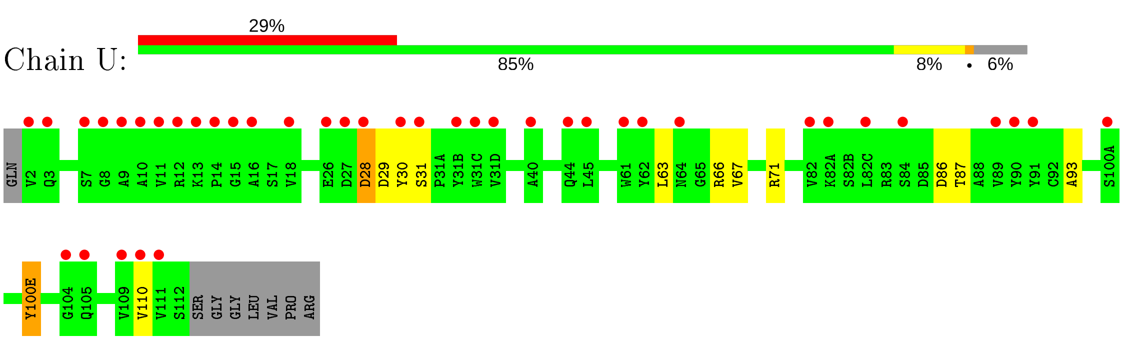
• Molecule 5: 3H109L Fab heavy chain



• Molecule 6: 3H109L Fab light chain

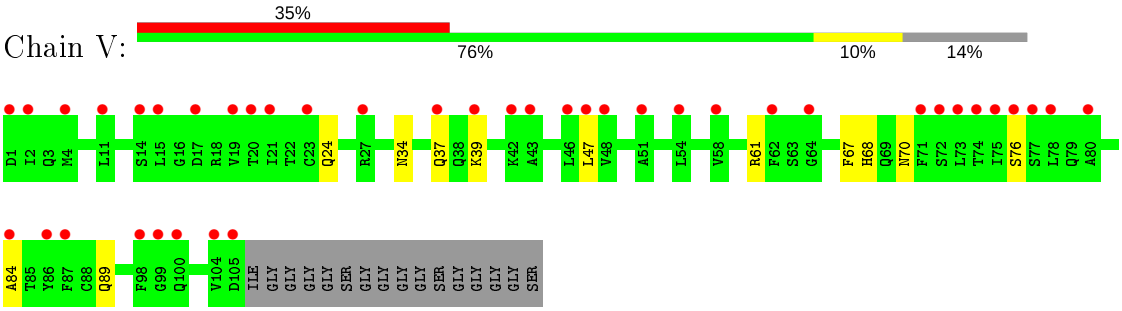


• Molecule 7: CH31 scFv heavy chain



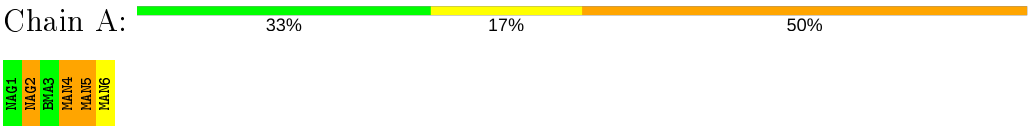
• Molecule 8: CH31 scFv light chain

Chain V:



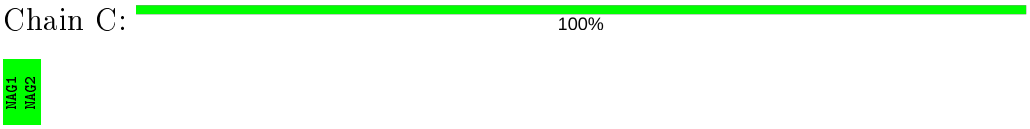
• Molecule 9: alpha-D-mannopyranose-(1-3)-alpha-D-mannopyranose-(1-6)-[alpha-D-mannopyranose-(1-3)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain A:



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

Chain C:

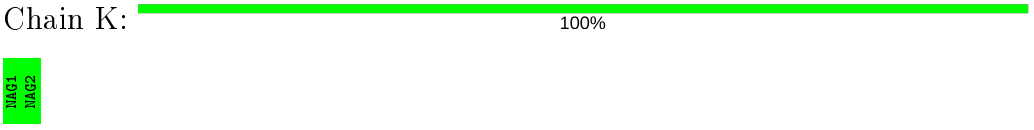


• Molecule 10: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain J:

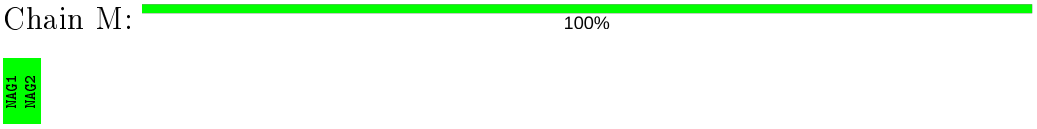
• Molecule 10: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain K:



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

Chain M:



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

Chain O:

• 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:

• 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 I:

• Molecule 13: 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

Chain N:

4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 63	Depositor
Cell constants a, b, c, α, β, γ	132.15Å 132.15Å 315.95Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	43.26 – 2.60 47.84 – 2.57	Depositor EDS
% Data completeness (in resolution range)	41.7 (43.26-2.60) 40.4 (47.84-2.57)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	1.54 (at 2.58Å)	Xtriage
Refinement program	PHENIX 1.14_3260	Depositor
R, R_{free}	0.248 , 0.293 0.248 , 0.293	Depositor DCC
R_{free} test set	1989 reflections (4.99%)	wwPDB-VP
Wilson B-factor (Å^2)	31.5	Xtriage
Anisotropy	0.107	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.28 , 26.7	EDS
L-test for twinning[2]	$< \|L\| > = 0.50$, $< L^2 > = 0.33$	Xtriage
Estimated twinning fraction	0.048 for h,-h-k,-l	Xtriage
F_o,F_c correlation	0.81	EDS
Total number of atoms	11874	wwPDB-VP
Average B, all atoms (Å^2)	55.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.01% of the height of the origin peak. No significant pseudotranslation is detected.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: BMA, NAG, 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	0.25	0/1027	0.39	0/1393
2	D	0.25	0/1021	0.48	0/1390
3	E	0.25	0/860	0.42	0/1175
4	G	0.27	0/3467	0.47	0/4702
5	H	0.26	0/1758	0.46	0/2397
6	L	0.25	0/1641	0.45	0/2239
7	U	0.24	0/1059	0.43	0/1447
8	V	0.25	0/774	0.46	0/1045
All	All	0.25	0/11607	0.45	0/15788

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	B	1009	0	994	16	0
2	D	994	0	953	20	0
3	E	836	0	785	8	0
4	G	3397	0	3349	73	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	27	0
7	U	1027	0	973	4	0
8	V	761	0	735	7	0
9	A	72	0	61	4	0
10	C	28	0	25	0	0
10	J	28	0	25	2	0
10	K	28	0	25	0	0
10	M	28	0	25	0	0
10	O	28	0	25	0	0
11	F	61	0	52	0	0
12	I	50	0	43	1	0
13	N	116	0	97	2	0
14	B	28	0	26	0	0
14	G	70	0	65	0	0
All	All	11874	0	11491	176	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 8.

All (176) 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)
5:H:59:TYR:HB2	5:H:64:LYS:HD2	1.56	0.87
4:G:230:ASP:HB3	4:G:233:PHE:HB2	1.62	0.81
4:G:475:MET:SD	4:G:478:ASN:ND2	2.55	0.78
4:G:358:ILE:HG22	4:G:396:ILE:HA	1.68	0.73
4:G:259:LEU:HD23	4:G:449:ILE:HG21	1.70	0.72
1:B:605:CYS:HA	4:G:37:THR:HG22	1.72	0.72
4:G:271:MET:HB3	4:G:273:ARG:HE	1.56	0.71
4:G:292:VAL:HG13	4:G:449:ILE:HG13	1.74	0.69
5:H:35:SER:OG	5:H:47:TRP:NE1	2.26	0.69
4:G:360:ARG:HB3	4:G:467:THR:HG22	1.74	0.68
5:H:157:LEU:HD21	5:H:180:VAL:HG11	1.75	0.67
5:H:100:ARG:NH2	13:N:4:MAN:O6	2.27	0.67
4:G:69:TRP:HE1	4:G:108:ILE:HG23	1.59	0.67
4:G:299:PRO:HG2	4:G:327:ARG:HB2	1.76	0.66
4:G:257:THR:HG22	4:G:258:GLN:HG3	1.77	0.66
5:H:35:SER:HG	5:H:47:TRP:HE1	1.41	0.66
4:G:360:ARG:HH11	4:G:467:THR:HG21	1.62	0.64
6:L:50:ASN:O	6:L:52:GLN:N	2.32	0.63
8:V:67:PHE:HZ	10:J:2:NAG:H4	1.63	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:96:LEU:HG	2:D:97:LEU:HG	1.79	0.62
8:V:24:GLN:NE2	8:V:70:ASN:OD1	2.33	0.62
1:B:520:LEU:HB3	1:B:524:GLY:HA3	1.81	0.61
5:H:63:LEU:O	5:H:65:SER:N	2.33	0.61
5:H:92:CYS:O	5:H:102:GLY:N	2.33	0.61
6:L:54:ARG:NH2	6:L:62:PHE:O	2.32	0.60
2:D:50:TRP:HH2	9:A:4:MAN:H62	1.66	0.59
4:G:335:LYS:HB3	4:G:412:ASP:HB3	1.83	0.59
4:G:476:ARG:HA	4:G:479:TRP:CD1	2.39	0.58
7:U:66:ARG:NH2	7:U:86:ASP:OD2	2.37	0.58
5:H:117:PRO:HB3	5:H:143:TYR:HB3	1.86	0.58
5:H:99:LYS:HE2	5:H:100(A):ILE:HD11	1.85	0.57
3:E:37:GLN:HG3	3:E:84:THR:HG21	1.86	0.57
1:B:629:LEU:HD23	4:G:44:VAL:HG23	1.86	0.57
5:H:68:ILE:HG23	5:H:81:LYS:HB2	1.86	0.57
4:G:42:VAL:HG23	4:G:44:VAL:HG12	1.87	0.57
4:G:201:ILE:HD11	4:G:435:TYR:HB2	1.87	0.57
4:G:324:GLY:O	6:L:94:ARG:NH1	2.34	0.56
1:B:585:ARG:NH2	4:G:491:ILE:O	2.36	0.56
9:A:2:NAG:H83	9:A:2:NAG:H3	1.88	0.55
4:G:66:HIS:ND1	4:G:66:HIS:N	2.55	0.55
5:H:195:ASN:ND2	5:H:206:ASP:OD2	2.40	0.54
1:B:519:PHE:HZ	1:B:542:ARG:HH22	1.55	0.54
4:G:350:ARG:NH2	4:G:357:THR:O	2.35	0.54
4:G:286:VAL:HG13	4:G:452:LEU:HB3	1.90	0.54
2:D:35:ASN:ND2	2:D:100(D):TRP:O	2.38	0.53
6:L:8:VAL:HG11	6:L:103:ARG:HE	1.73	0.53
7:U:87:THR:HG23	7:U:110:VAL:HA	1.90	0.53
6:L:35:TRP:CE2	6:L:73:LEU:HB2	2.43	0.53
4:G:102:GLU:OE2	4:G:476:ARG:NH1	2.42	0.52
6:L:34:GLN:HG3	6:L:49:TYR:HA	1.91	0.52
8:V:39:LYS:HG2	8:V:84:ALA:HB2	1.90	0.52
2:D:66:ARG:NH1	2:D:86:ASP:OD2	2.42	0.52
5:H:136:LEU:HD13	5:H:209:VAL:HG21	1.92	0.52
8:V:37:GLN:HB2	8:V:47:LEU:HD11	1.90	0.52
6:L:109:GLN:HB2	6:L:141:TYR:CE1	2.45	0.52
5:H:165:PRO:HG2	6:L:163:THR:HG21	1.91	0.51
6:L:37:GLN:HB2	6:L:47:LEU:HD11	1.92	0.51
2:D:50:TRP:CH2	9:A:4:MAN:H62	2.45	0.51
6:L:124:GLU:OE1	6:L:124:GLU:N	2.43	0.51
4:G:257:THR:HG21	4:G:370:GLU:O	2.11	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:L:28:LEU:HB2	6:L:94:ARG:HB2	1.93	0.51
6:L:14:ALA:HB3	6:L:17:GLU:HG3	1.92	0.51
8:V:67:PHE:CZ	10:J:2:NAG:H4	2.46	0.50
6:L:119:PHE:HB2	6:L:134:VAL:HG13	1.92	0.50
5:H:142:ASP:OD1	5:H:142:ASP:N	2.43	0.50
3:E:37:GLN:HB2	3:E:47:ILE:HD11	1.94	0.50
2:D:48:MET:HG2	2:D:63:PHE:CD2	2.47	0.50
5:H:4:LEU:HG	5:H:24:VAL:HG23	1.93	0.50
4:G:69:TRP:HE1	4:G:108:ILE:HD12	1.76	0.49
4:G:123:THR:N	4:G:124:PRO:HD2	2.28	0.49
3:E:83:GLU:HG3	3:E:106:VAL:HG12	1.95	0.49
4:G:112:TRP:CE2	4:G:426:MET:HG2	2.48	0.49
1:B:571:TRP:CZ2	4:G:111:LEU:HD11	2.48	0.48
4:G:439:ILE:HB	4:G:443:ILE:HD11	1.95	0.48
4:G:299:PRO:HA	4:G:442:VAL:HG13	1.94	0.48
4:G:456:ARG:NH1	4:G:466:GLU:OE2	2.45	0.48
1:B:569:THR:HG23	1:B:572:GLY:H	1.78	0.48
4:G:69:TRP:NE1	4:G:108:ILE:HG23	2.27	0.48
4:G:369:LEU:O	4:G:373:THR:OG1	2.29	0.48
2:D:32:TYR:CD2	2:D:94:LYS:HE3	2.49	0.48
6:L:33:VAL:HG12	6:L:51:ASN:OD1	2.14	0.48
6:L:19:ALA:HB3	6:L:75:ILE:HB	1.95	0.48
4:G:122:LEU:HD11	4:G:203:GLN:HB2	1.95	0.48
5:H:103:LYS:HD2	5:H:103:LYS:N	2.29	0.47
4:G:343:GLY:O	4:G:346:VAL:HG12	2.14	0.47
2:D:36:TRP:CZ3	2:D:92:CYS:HB3	2.50	0.47
7:U:93:ALA:HB1	7:U:100(E):TYR:HB3	1.96	0.47
13:N:1:NAG:H3	13:N:1:NAG:H83	1.96	0.47
6:L:197:THR:HA	6:L:202:THR:HA	1.98	0.46
4:G:270:VAL:HG11	4:G:344:LYS:HB3	1.97	0.46
6:L:83:GLU:OE1	6:L:167:LYS:NZ	2.48	0.46
2:D:13:LYS:HD2	2:D:13:LYS:HA	1.68	0.46
2:D:38:ARG:HD2	2:D:46:GLU:HB3	1.97	0.46
1:B:610:TRP:CD2	4:G:498:PRO:HB3	2.52	0.45
4:G:105:HIS:CD2	4:G:476:ARG:HH21	2.34	0.45
5:H:6:GLU:N	5:H:6:GLU:OE1	2.47	0.45
3:E:27(C):CYS:HA	3:E:28:CYS:HA	1.66	0.45
4:G:358:ILE:HG13	4:G:465:THR:HG22	1.98	0.45
4:G:333:VAL:HG21	4:G:390:LEU:HD21	1.97	0.45
4:G:94:ASN:HB3	4:G:97:LYS:HG2	1.98	0.45
4:G:54:CYS:SG	4:G:55:ALA:N	2.90	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:638:TYR:O	1:B:642:ILE:HG13	2.16	0.45
4:G:104:MET:O	4:G:108:ILE:HG12	2.17	0.45
4:G:206:PRO:HD2	4:G:207:LYS:HZ2	1.82	0.45
1:B:606:THR:OG1	4:G:36:VAL:O	2.33	0.44
3:E:7:SER:HB2	3:E:22:SER:H	1.83	0.44
2:D:28:ARG:HG2	2:D:72(H):PHE:O	2.17	0.44
1:B:574:LYS:HE3	4:G:107:ASP:OD2	2.18	0.44
6:L:59:PRO:HB2	6:L:61:ARG:HG2	1.99	0.44
6:L:39:ARG:HG2	6:L:84:ALA:HB2	2.00	0.44
2:D:30:ASN:HA	2:D:52(A):PRO:HB2	2.00	0.44
4:G:122:LEU:HB3	4:G:125:LEU:HD12	2.00	0.44
2:D:59:LEU:HD11	2:D:64:GLN:HA	1.98	0.44
4:G:164:GLU:OE2	4:G:308:ARG:NH1	2.51	0.44
9:A:4:MAN:H3	9:A:5:MAN:H2	1.68	0.43
5:H:27:GLY:O	5:H:76:ASN:ND2	2.51	0.43
6:L:83:GLU:HG3	6:L:106:VAL:HG23	1.99	0.43
6:L:35:TRP:CD2	6:L:73:LEU:HB2	2.53	0.43
4:G:256:SER:OG	4:G:257:THR:N	2.51	0.43
4:G:391:PHE:CD2	4:G:470:PRO:HG3	2.53	0.43
6:L:143:GLY:HA3	6:L:173:TYR:CG	2.53	0.43
4:G:385:CYS:HA	4:G:418:CYS:HA	2.00	0.43
4:G:86:LEU:HB3	4:G:89:VAL:HG21	2.01	0.43
5:H:191:THR:HB	5:H:208:LYS:HE2	2.00	0.43
4:G:234:ASN:OD1	4:G:235:GLY:N	2.52	0.43
1:B:618:ASN:O	1:B:621:GLU:HG2	2.18	0.43
3:E:42:ARG:HB3	3:E:43:ALA:H	1.63	0.43
4:G:206:PRO:HD2	4:G:207:LYS:NZ	2.33	0.43
4:G:260:LEU:HD12	4:G:451:GLY:HA3	2.00	0.43
4:G:206:PRO:HG3	4:G:318:TYR:CE2	2.54	0.43
4:G:257:THR:HB	4:G:375:SER:H	1.83	0.43
4:G:261:LEU:HD13	12:I:1:NAG:H82	2.00	0.43
5:H:142:ASP:HB3	5:H:173:LEU:HD22	2.01	0.42
2:D:6:GLN:H	2:D:105:GLN:NE2	2.17	0.42
4:G:134:VAL:HG21	4:G:154:LEU:HD12	2.01	0.42
4:G:359:ILE:HD12	4:G:468:PHE:HE2	1.84	0.42
4:G:84:ILE:HB	4:G:244:THR:HG23	2.01	0.42
3:E:78:LEU:HD13	3:E:106:VAL:HG23	2.02	0.42
5:H:144:PHE:HA	5:H:145:PRO:HA	1.81	0.42
4:G:66:HIS:O	4:G:208:VAL:HA	2.18	0.42
5:H:24:VAL:HG12	5:H:76:ASN:HB3	2.02	0.42
1:B:627:THR:OG1	1:B:630:GLN:HG3	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:L:36:TYR:CE1	6:L:46:LEU:HD13	2.54	0.42
2:D:36:TRP:HD1	2:D:69:MET:SD	2.42	0.42
2:D:19:LYS:HE3	2:D:79:TYR:HB3	2.02	0.42
4:G:69:TRP:NE1	4:G:108:ILE:HD12	2.33	0.42
5:H:24:VAL:CG1	5:H:76:ASN:HB3	2.50	0.42
4:G:95:MET:SD	4:G:235:GLY:HA2	2.60	0.41
5:H:100(A):ILE:HG21	5:H:100(E):VAL:HG22	2.02	0.41
6:L:47:LEU:O	6:L:48:ILE:HG13	2.20	0.41
8:V:61:ARG:HB2	8:V:76:SER:O	2.20	0.41
3:E:18:SER:HB3	3:E:76:SER:HA	2.03	0.41
4:G:216:HIS:ND1	4:G:248:THR:O	2.49	0.41
5:H:120:PHE:HE2	5:H:141:LYS:HE2	1.84	0.41
5:H:8:GLY:HA3	5:H:20:LEU:HD23	2.01	0.41
6:L:13:VAL:HG21	6:L:78:VAL:HG21	2.03	0.41
1:B:569:THR:OG1	1:B:570:VAL:N	2.53	0.41
2:D:11:THR:HG22	2:D:109:LEU:O	2.21	0.41
2:D:94:LYS:HB3	2:D:102:LEU:HB3	2.03	0.41
2:D:68:ASN:HB2	2:D:81:GLU:HB2	2.02	0.41
4:G:205:CYS:HB3	4:G:207:LYS:HD2	2.02	0.41
4:G:338:TRP:CE2	4:G:390:LEU:HD22	2.56	0.41
4:G:213:ILE:HA	4:G:213:ILE:HD13	1.89	0.41
5:H:91:TYR:CE2	6:L:43:ALA:HA	2.55	0.41
6:L:168:GLN:OE1	6:L:174:ALA:HB2	2.21	0.41
8:V:34:ASN:O	8:V:89:GLN:N	2.54	0.41
2:D:57:LYS:HE3	2:D:69:MET:HE2	2.03	0.41
1:B:523:LEU:O	4:G:86:LEU:HD22	2.21	0.41
4:G:360:ARG:NH1	4:G:467:THR:HG21	2.34	0.40
1:B:574:LYS:NZ	4:G:103:GLN:OE1	2.54	0.40
4:G:295:ASN:O	4:G:331:CYS:HA	2.22	0.40
5:H:47:TRP:CZ2	5:H:100(P):MET:HE1	2.55	0.40
7:U:63:LEU:HD13	7:U:67:VAL:HG21	2.04	0.40
5:H:100:ARG:HA	5:H:100:ARG:HD2	1.86	0.40
4:G:358:ILE:H	4:G:358:ILE:HG12	1.70	0.40
5:H:83:THR:OG1	5:H:84:ALA:N	2.54	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	123/153 (80%)	116 (94%)	7 (6%)	0	100	100
2	D	126/153 (82%)	115 (91%)	10 (8%)	1 (1%)	19	39
3	E	108/130 (83%)	89 (82%)	17 (16%)	2 (2%)	8	15
4	G	418/481 (87%)	383 (92%)	31 (7%)	4 (1%)	15	32
5	H	222/244 (91%)	200 (90%)	18 (8%)	4 (2%)	8	16
6	L	208/217 (96%)	194 (93%)	11 (5%)	3 (1%)	11	22
7	U	127/137 (93%)	120 (94%)	4 (3%)	3 (2%)	6	10
8	V	99/117 (85%)	91 (92%)	7 (7%)	1 (1%)	15	32
All	All	1431/1632 (88%)	1308 (91%)	105 (7%)	18 (1%)	12	24

All (18) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
6	L	51	ASN
7	U	31	SER
4	G	427	TRP
6	L	143	GLY
4	G	70	ALA
4	G	275	GLU
5	H	64	LYS
6	L	152	ASP
3	E	8	ALA
3	E	77	ASP
8	V	68	HIS
4	G	426	MET
5	H	117	PRO
7	U	28	ASP
7	U	100(E)	TYR
5	H	124	PRO
2	D	100(F)	PRO

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Mol	Chain	Res	Type
5	H	147	PRO

5.3.2 Protein sidechains

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
1	B	109/129 (84%)	107 (98%)	2 (2%)	59 80
2	D	107/115 (93%)	102 (95%)	5 (5%)	26 50
3	E	96/113 (85%)	92 (96%)	4 (4%)	30 55
4	G	385/427 (90%)	372 (97%)	13 (3%)	37 63
5	H	196/212 (92%)	190 (97%)	6 (3%)	40 66
6	L	174/181 (96%)	169 (97%)	5 (3%)	42 68
7	U	105/111 (95%)	101 (96%)	4 (4%)	33 59
8	V	84/88 (96%)	84 (100%)	0	100 100
All	All	1256/1376 (91%)	1217 (97%)	39 (3%)	40 66

All (39) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	548	ILE
1	B	569	THR
2	D	38	ARG
2	D	68	ASN
2	D	91	PHE
2	D	100(C)	THR
2	D	105	GLN
3	E	19	VAL
3	E	68	TYR
3	E	89	CYS
3	E	96	CYS
4	G	54	CYS
4	G	57	ASP
4	G	65	LYS

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Mol	Chain	Res	Type
4	G	66	HIS
4	G	154	LEU
4	G	207	LYS
4	G	270	VAL
4	G	286	VAL
4	G	292	VAL
4	G	346	VAL
4	G	432	GLN
4	G	444	ARG
4	G	447	SER
5	H	100(F)	SER
5	H	100(P)	MET
5	H	142	ASP
5	H	162	HIS
5	H	203	THR
5	H	207	LYS
6	L	25	ARG
6	L	54	ARG
6	L	97	SER
6	L	197	THR
6	L	205	LYS
7	U	28	ASP
7	U	29	ASP
7	U	30	TYR
7	U	71	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
8	V	24	GLN

5.3.3 RNA (i)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates (i)

35 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	0.38	0	17,19,21	0.47	0
9	NAG	A	2	9	14,14,15	0.41	0	17,19,21	1.32	1 (5%)
9	BMA	A	3	9	11,11,12	0.74	0	15,15,17	0.86	0
9	MAN	A	4	9	11,11,12	0.71	0	15,15,17	1.07	2 (13%)
9	MAN	A	5	9	11,11,12	1.73	1 (9%)	15,15,17	1.69	2 (13%)
9	MAN	A	6	9	11,11,12	0.72	0	15,15,17	0.99	2 (13%)
10	NAG	C	1	10,4	14,14,15	0.22	0	17,19,21	0.44	0
10	NAG	C	2	10	14,14,15	0.28	0	17,19,21	0.47	0
11	NAG	F	1	11,4	14,14,15	0.30	0	17,19,21	0.46	0
11	NAG	F	2	11	14,14,15	0.16	0	17,19,21	0.41	0
11	BMA	F	3	11	11,11,12	0.72	0	15,15,17	0.75	0
11	MAN	F	4	11	11,11,12	0.75	0	15,15,17	1.05	2 (13%)
11	MAN	F	5	11	11,11,12	0.78	1 (9%)	15,15,17	1.03	2 (13%)
12	NAG	I	1	12,4	14,14,15	0.33	0	17,19,21	0.47	0
12	NAG	I	2	12	14,14,15	0.23	0	17,19,21	0.58	0
12	BMA	I	3	12	11,11,12	0.72	0	15,15,17	0.76	0
12	MAN	I	4	12	11,11,12	0.78	1 (9%)	15,15,17	1.12	2 (13%)
10	NAG	J	1	10,4	14,14,15	0.32	0	17,19,21	0.38	0
10	NAG	J	2	10	14,14,15	0.35	0	17,19,21	0.44	0
10	NAG	K	1	10,4	14,14,15	0.34	0	17,19,21	0.38	0
10	NAG	K	2	10	14,14,15	0.20	0	17,19,21	0.45	0
10	NAG	M	1	10,4	14,14,15	0.30	0	17,19,21	0.46	0
10	NAG	M	2	10	14,14,15	0.27	0	17,19,21	0.51	0
13	NAG	N	1	13,4	14,14,15	0.28	0	17,19,21	1.51	2 (11%)
13	MAN	N	10	13	11,11,12	1.11	2 (18%)	15,15,17	1.58	4 (26%)
13	NAG	N	2	13	14,14,15	0.53	0	17,19,21	0.46	0
13	BMA	N	3	13	11,11,12	1.02	0	15,15,17	1.08	1 (6%)
13	MAN	N	4	13	11,11,12	0.82	1 (9%)	15,15,17	1.34	1 (6%)
13	MAN	N	5	13	11,11,12	0.76	0	15,15,17	1.16	1 (6%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\|>2$	Counts	RMSZ	$\#\|Z\|>2$
13	MAN	N	6	13	11,11,12	0.91	0	15,15,17	0.92	1 (6%)
13	MAN	N	7	13	11,11,12	0.74	0	15,15,17	0.97	2 (13%)
13	MAN	N	8	13	11,11,12	0.70	0	15,15,17	1.01	1 (6%)
13	MAN	N	9	13	11,11,12	0.89	1 (9%)	15,15,17	1.26	2 (13%)
10	NAG	O	1	10,4	14,14,15	0.92	1 (7%)	17,19,21	1.48	1 (5%)
10	NAG	O	2	10	14,14,15	0.26	0	17,19,21	0.37	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	-	5/6/23/26	0/1/1/1
9	BMA	A	3	9	-	2/2/19/22	0/1/1/1
9	MAN	A	4	9	-	0/2/19/22	0/1/1/1
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	-	0/6/23/26	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
12	NAG	I	1	12,4	-	2/6/23/26	0/1/1/1
12	NAG	I	2	12	-	2/6/23/26	0/1/1/1
12	BMA	I	3	12	-	2/2/19/22	0/1/1/1
12	MAN	I	4	12	-	2/2/19/22	0/1/1/1
10	NAG	J	1	10,4	-	2/6/23/26	0/1/1/1
10	NAG	J	2	10	-	2/6/23/26	0/1/1/1
10	NAG	K	1	10,4	-	2/6/23/26	0/1/1/1
10	NAG	K	2	10	-	0/6/23/26	0/1/1/1
10	NAG	M	1	10,4	-	1/6/23/26	0/1/1/1
10	NAG	M	2	10	-	0/6/23/26	0/1/1/1
13	NAG	N	1	13,4	-	5/6/23/26	0/1/1/1
13	MAN	N	10	13	-	0/2/19/22	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
13	NAG	N	2	13	-	2/6/23/26	0/1/1/1
13	BMA	N	3	13	-	0/2/19/22	0/1/1/1
13	MAN	N	4	13	-	2/2/19/22	0/1/1/1
13	MAN	N	5	13	-	0/2/19/22	0/1/1/1
13	MAN	N	6	13	-	0/2/19/22	0/1/1/1
13	MAN	N	7	13	-	0/2/19/22	0/1/1/1
13	MAN	N	8	13	-	0/2/19/22	0/1/1/1
13	MAN	N	9	13	-	0/2/19/22	0/1/1/1
10	NAG	O	1	10,4	-	4/6/23/26	0/1/1/1
10	NAG	O	2	10	-	2/6/23/26	0/1/1/1

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
9	A	5	MAN	C1-C2	5.07	1.63	1.52
10	O	1	NAG	O5-C1	3.32	1.49	1.43
13	N	9	MAN	C1-C2	2.49	1.57	1.52
13	N	10	MAN	C1-C2	2.42	1.57	1.52
12	I	4	MAN	C1-C2	2.27	1.57	1.52
13	N	4	MAN	C1-C2	2.11	1.57	1.52
13	N	10	MAN	C2-C3	2.09	1.55	1.52
11	F	5	MAN	C1-C2	2.03	1.56	1.52

All (29) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
10	O	1	NAG	C1-O5-C5	5.73	119.96	112.19
13	N	1	NAG	C2-N2-C7	4.71	129.61	122.90
9	A	5	MAN	C1-C2-C3	4.56	115.27	109.67
9	A	2	NAG	C2-N2-C7	4.38	129.14	122.90
13	N	4	MAN	C1-O5-C5	3.84	117.39	112.19
13	N	10	MAN	C1-C2-C3	3.54	114.01	109.67
13	N	9	MAN	C1-O5-C5	3.39	116.79	112.19
9	A	5	MAN	C1-O5-C5	3.29	116.65	112.19
13	N	5	MAN	C1-O5-C5	3.03	116.30	112.19
12	I	4	MAN	C1-O5-C5	2.63	115.75	112.19
13	N	1	NAG	C1-C2-N2	2.61	114.95	110.49
9	A	4	MAN	C1-O5-C5	2.51	115.59	112.19
9	A	6	MAN	O2-C2-C3	-2.46	105.20	110.14
9	A	6	MAN	C1-O5-C5	2.39	115.43	112.19
13	N	9	MAN	O2-C2-C3	-2.33	105.48	110.14

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
13	N	8	MAN	C1-O5-C5	2.31	115.32	112.19
13	N	10	MAN	O2-C2-C3	-2.30	105.53	110.14
11	F	5	MAN	C1-O5-C5	2.29	115.29	112.19
13	N	7	MAN	O2-C2-C3	-2.29	105.55	110.14
13	N	10	MAN	C1-O5-C5	2.27	115.27	112.19
13	N	10	MAN	O5-C1-C2	2.24	114.22	110.77
13	N	3	BMA	C1-C2-C3	2.22	112.40	109.67
13	N	6	MAN	O2-C2-C3	-2.21	105.71	110.14
11	F	4	MAN	O2-C2-C3	-2.13	105.88	110.14
12	I	4	MAN	O2-C2-C3	-2.12	105.89	110.14
9	A	4	MAN	O2-C2-C3	-2.11	105.92	110.14
11	F	5	MAN	O2-C2-C3	-2.10	105.93	110.14
13	N	7	MAN	C1-O5-C5	2.05	114.97	112.19
11	F	4	MAN	C1-O5-C5	2.02	114.94	112.19

There are no chirality outliers.

All (47) torsion outliers are listed below:

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

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Mol	Chain	Res	Type	Atoms
10	O	2	NAG	C4-C5-C6-O6
10	J	1	NAG	O5-C5-C6-O6
13	N	1	NAG	O5-C5-C6-O6
10	K	1	NAG	C4-C5-C6-O6
11	F	4	MAN	C4-C5-C6-O6
13	N	4	MAN	O5-C5-C6-O6
12	I	3	BMA	C4-C5-C6-O6
13	N	1	NAG	C4-C5-C6-O6
10	J	1	NAG	C4-C5-C6-O6
11	F	3	BMA	O5-C5-C6-O6
13	N	4	MAN	C4-C5-C6-O6
11	F	5	MAN	C4-C5-C6-O6
10	C	2	NAG	C4-C5-C6-O6
9	A	5	MAN	O5-C5-C6-O6
11	F	5	MAN	O5-C5-C6-O6
10	C	2	NAG	O5-C5-C6-O6
10	M	1	NAG	O5-C5-C6-O6
10	J	2	NAG	C4-C5-C6-O6
10	J	2	NAG	O5-C5-C6-O6
13	N	2	NAG	C4-C5-C6-O6
13	N	2	NAG	O5-C5-C6-O6
13	N	1	NAG	C3-C2-N2-C7
9	A	2	NAG	C3-C2-N2-C7
9	A	5	MAN	C4-C5-C6-O6

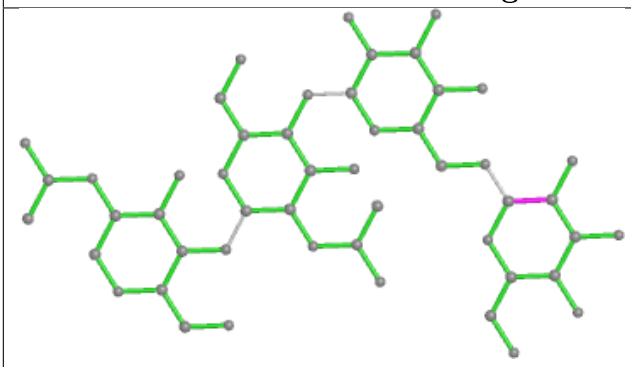
There are no ring outliers.

7 monomers are involved in 9 short contacts:

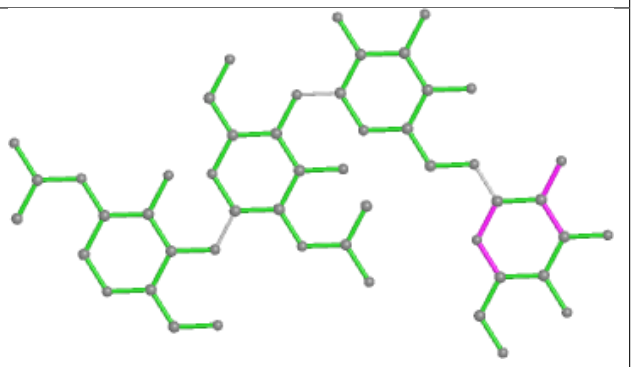
Mol	Chain	Res	Type	Clashes	Symm-Clashes
10	J	2	NAG	2	0
13	N	1	NAG	1	0
9	A	2	NAG	1	0
9	A	5	MAN	1	0
9	A	4	MAN	3	0
13	N	4	MAN	1	0
12	I	1	NAG	1	0

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

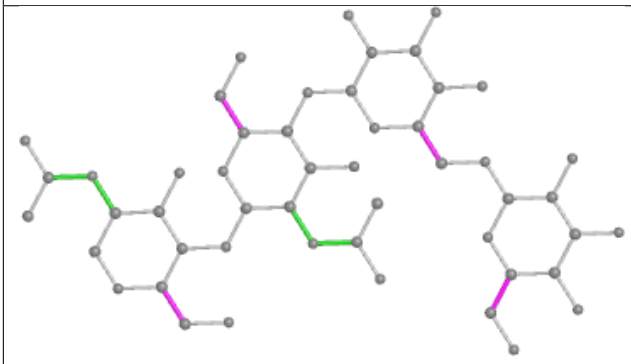
Oligosaccharide Chain I



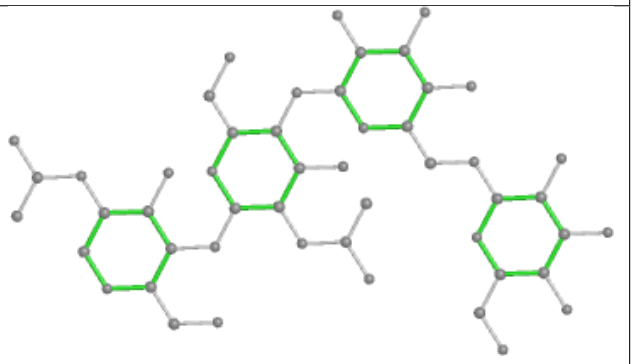
Bond lengths



Bond angles

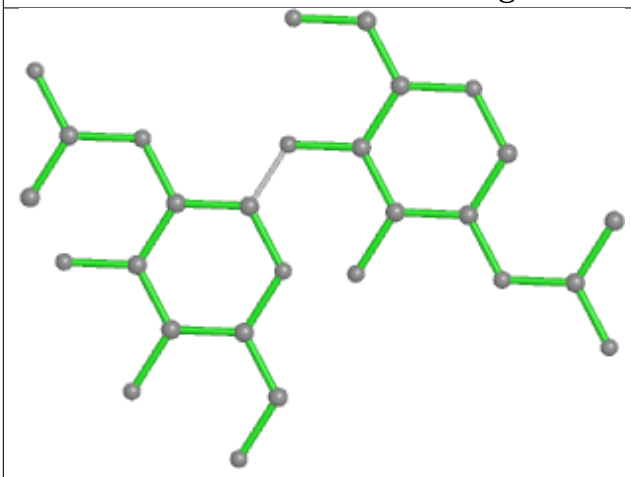


Torsions

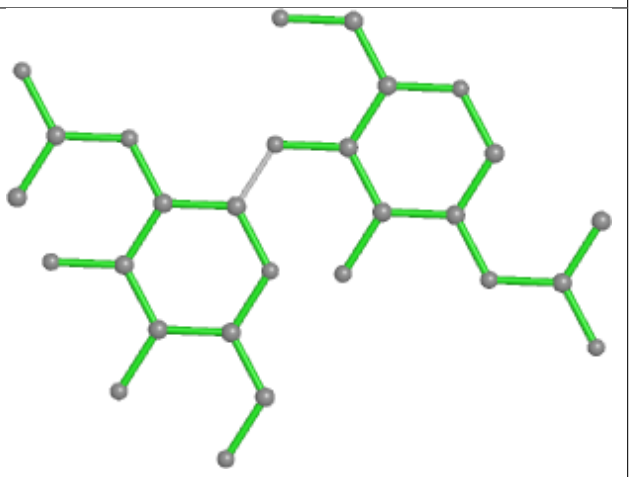


Rings

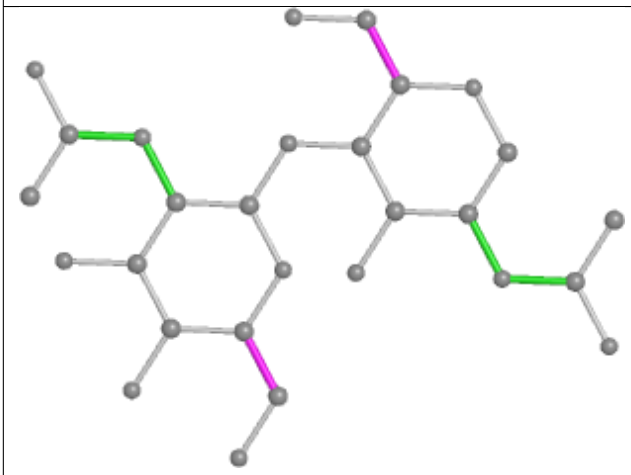
Oligosaccharide Chain J



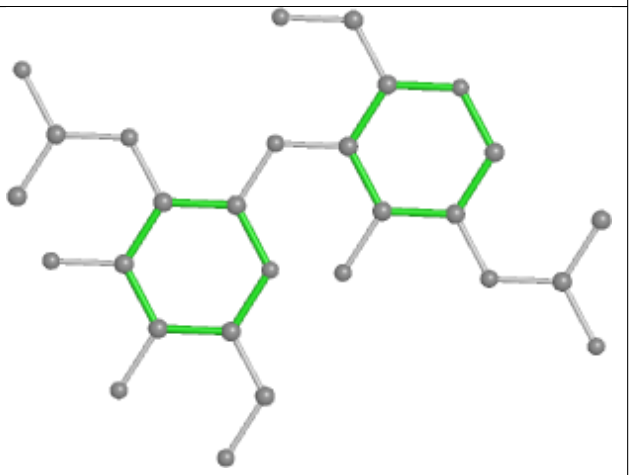
Bond lengths



Bond angles

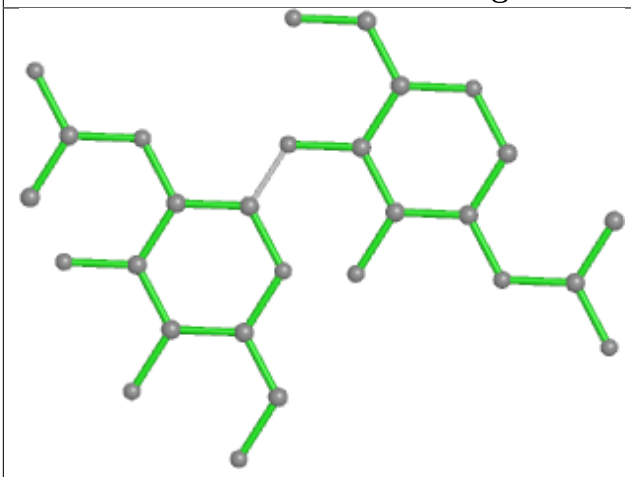


Torsions

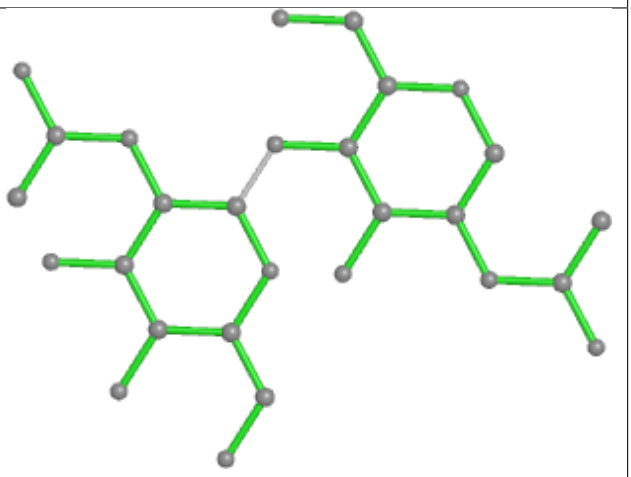


Rings

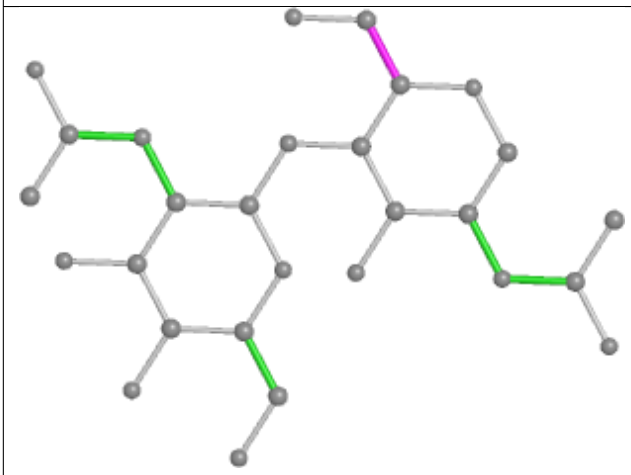
Oligosaccharide Chain M



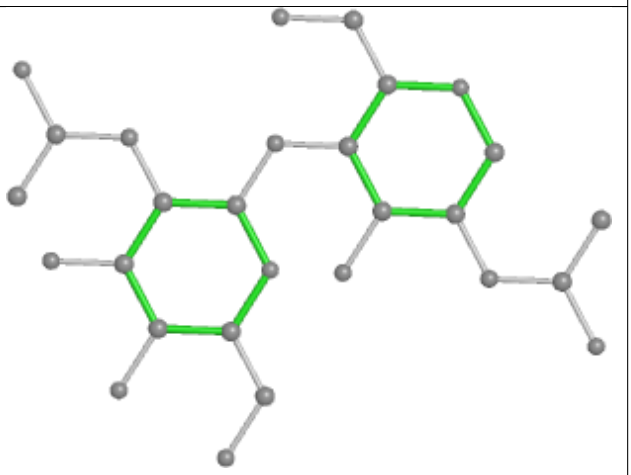
Bond lengths



Bond angles

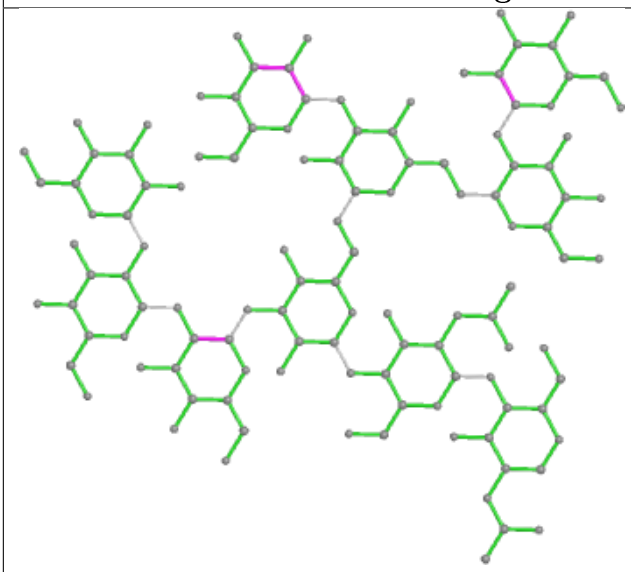


Torsions

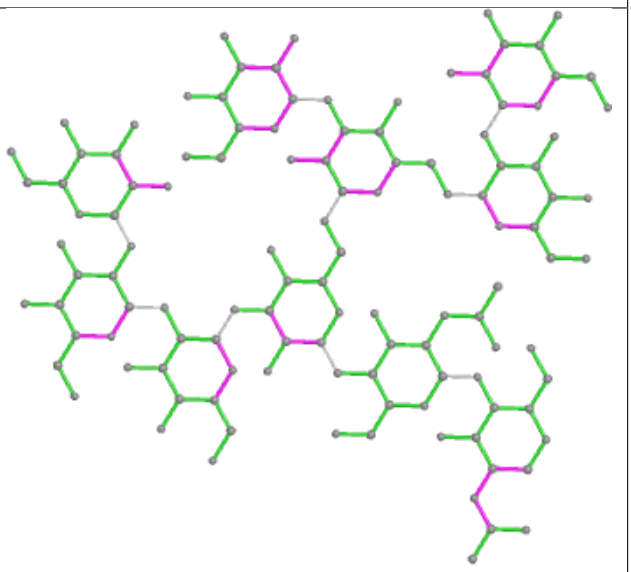


Rings

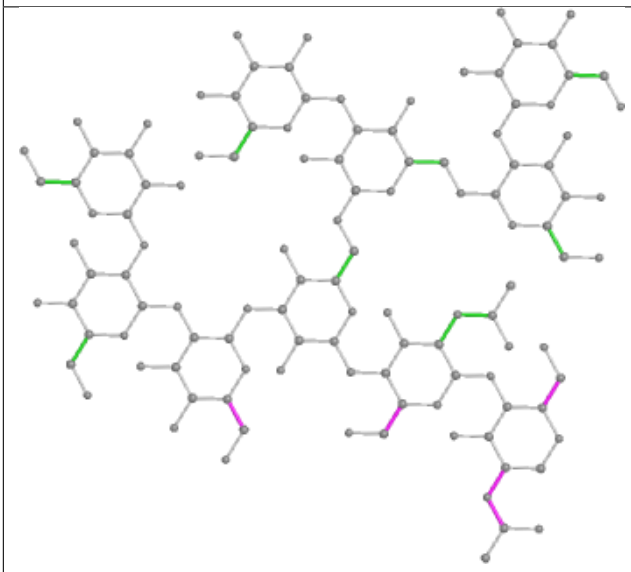
Oligosaccharide Chain N



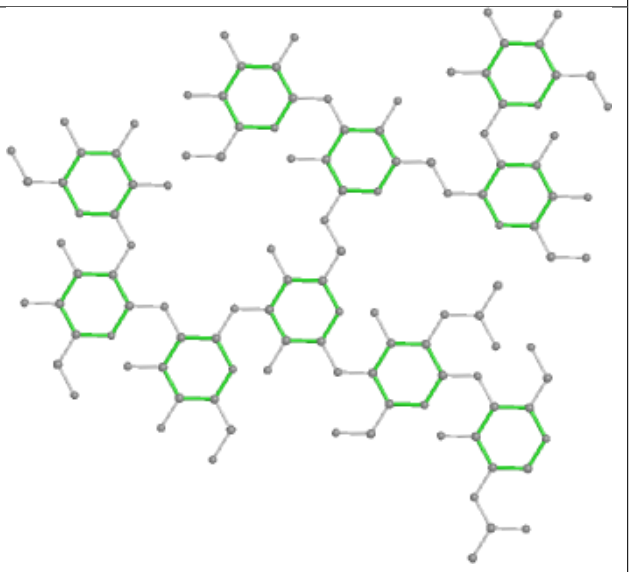
Bond lengths



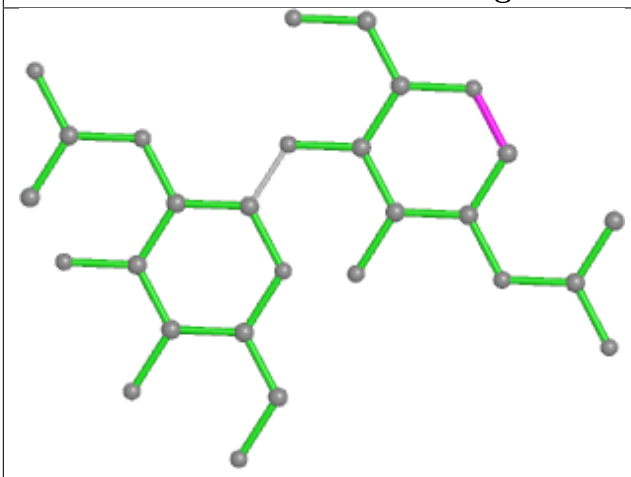
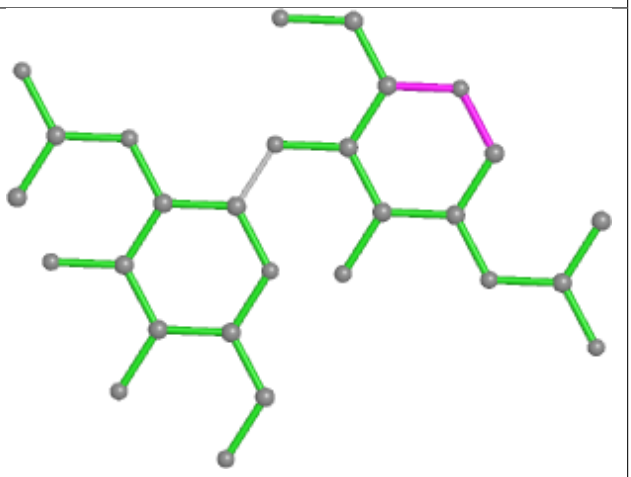
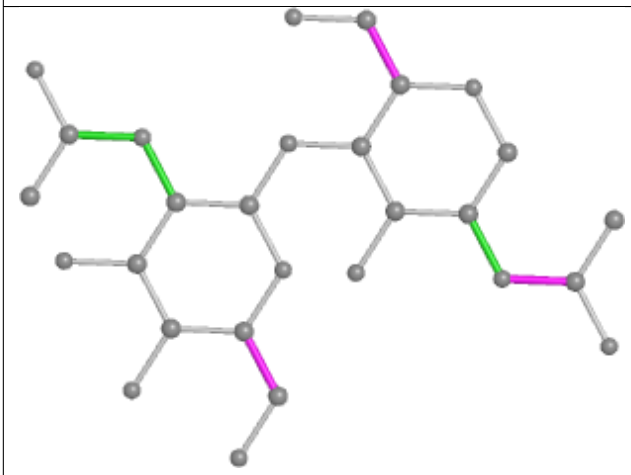
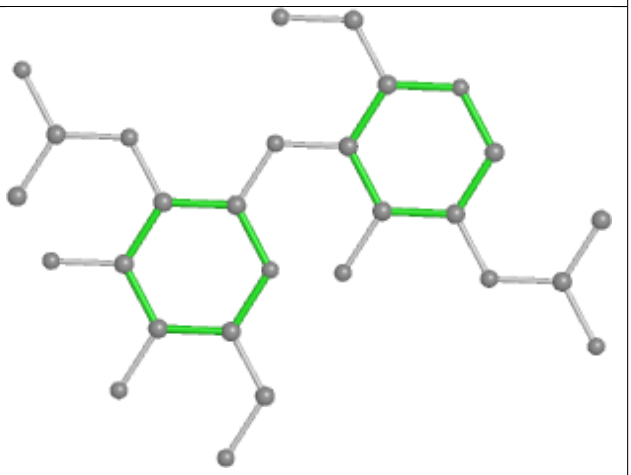
Bond angles

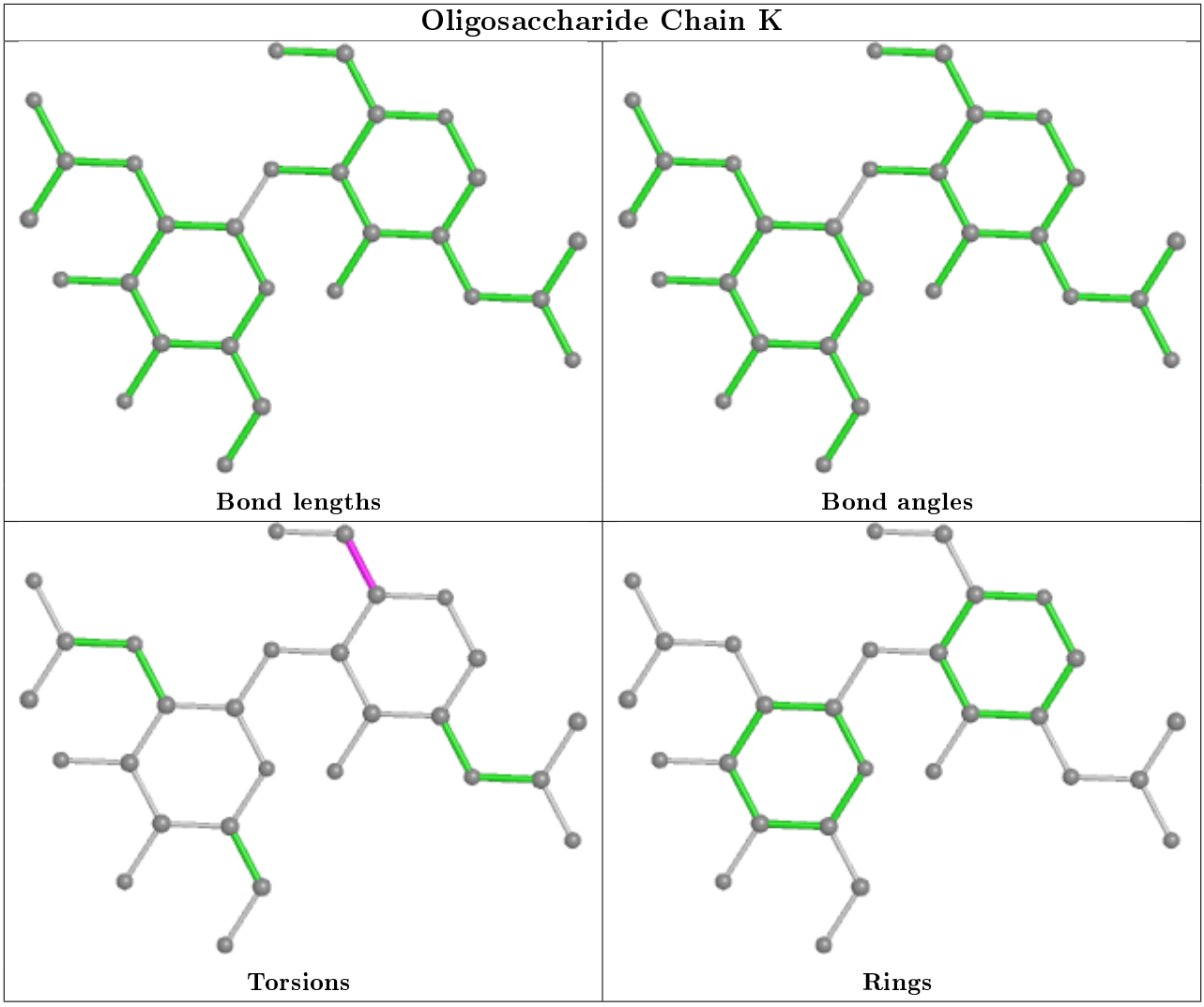


Torsions



Rings

Oligosaccharide Chain O	
	
Bond lengths	Bond angles
	
Torsions	Rings



5.6 Ligand geometry (i)

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	702	1	14,14,15	0.32	0	17,19,21	0.42	0
14	NAG	G	616	4	14,14,15	0.28	0	17,19,21	0.49	0
14	NAG	B	701	1	14,14,15	0.26	0	17,19,21	0.57	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\|>2$	Counts	RMSZ	$\#\|Z\|>2$
14	NAG	G	615	4	14,14,15	0.72	1 (7%)	17,19,21	0.58	0
14	NAG	G	638	4	14,14,15	0.26	0	17,19,21	0.51	0
14	NAG	G	637	4	14,14,15	0.26	0	17,19,21	0.50	0
14	NAG	G	614	4	14,14,15	0.28	0	17,19,21	0.49	0

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
14	NAG	B	702	1	-	2/6/23/26	0/1/1/1
14	NAG	G	616	4	-	2/6/23/26	0/1/1/1
14	NAG	B	701	1	-	2/6/23/26	0/1/1/1
14	NAG	G	615	4	-	2/6/23/26	0/1/1/1
14	NAG	G	638	4	-	2/6/23/26	0/1/1/1
14	NAG	G	637	4	-	2/6/23/26	0/1/1/1
14	NAG	G	614	4	-	2/6/23/26	0/1/1/1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
14	G	615	NAG	C1-C2	2.48	1.56	1.52

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	638	NAG	O5-C5-C6-O6
14	G	614	NAG	O5-C5-C6-O6
14	G	615	NAG	O5-C5-C6-O6
14	G	616	NAG	O5-C5-C6-O6
14	G	637	NAG	O5-C5-C6-O6
14	G	638	NAG	C4-C5-C6-O6
14	G	616	NAG	C4-C5-C6-O6
14	G	637	NAG	C4-C5-C6-O6
14	B	701	NAG	O5-C5-C6-O6
14	G	615	NAG	C4-C5-C6-O6

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Mol	Chain	Res	Type	Atoms
14	B	702	NAG	O5-C5-C6-O6
14	G	614	NAG	C4-C5-C6-O6
14	B	701	NAG	C4-C5-C6-O6
14	B	702	NAG	C4-C5-C6-O6

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers (i)

There are no such residues in this entry.

5.8 Polymer linkage issues (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, 95^{th} 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	127/153 (83%)	-0.22	0 100 100	8, 31, 61, 80	0
2	D	128/153 (83%)	1.32	37 (28%) 0 0	38, 83, 116, 127	0
3	E	110/130 (84%)	0.74	12 (10%) 5 3	42, 72, 104, 116	0
4	G	430/481 (89%)	-0.18	7 (1%) 72 68	8, 26, 71, 109	0
5	H	226/244 (92%)	0.33	22 (9%) 7 5	23, 61, 93, 105	0
6	L	210/217 (96%)	-0.20	4 (1%) 66 62	22, 46, 68, 95	0
7	U	129/137 (94%)	1.24	40 (31%) 0 0	34, 85, 116, 132	0
8	V	101/117 (86%)	1.94	41 (40%) 0 0	64, 101, 118, 123	0
All	All	1461/1632 (89%)	0.37	163 (11%) 5 3	8, 52, 107, 132	0

All (163) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	D	18	VAL	8.2
2	D	82	ILE	8.1
2	D	110	THR	7.6
8	V	2	ILE	7.6
8	V	21	ILE	7.3
8	V	19	VAL	6.6
2	D	41	ALA	6.6
2	D	82(C)	LEU	6.3
8	V	80	ALA	6.3
2	D	10	THR	6.2
3	E	106	VAL	6.1
2	D	109	LEU	6.0
2	D	13	LYS	5.9
8	V	20	THR	5.8
7	U	40	ALA	5.7
7	U	91	TYR	5.7

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Mol	Chain	Res	Type	RSRZ
8	V	99	GLY	5.4
7	U	18	VAL	5.3
5	H	187	LEU	5.3
3	E	68	TYR	5.0
7	U	109	VAL	4.9
8	V	105	ASP	4.7
8	V	73	LEU	4.6
8	V	47	LEU	4.6
7	U	26	GLU	4.5
2	D	91	PHE	4.5
7	U	12	ARG	4.5
8	V	76	SER	4.4
5	H	208	LYS	4.4
8	V	75	ILE	4.4
5	H	191	THR	4.4
5	H	209	VAL	4.3
2	D	85	ASP	4.3
7	U	11	VAL	4.3
8	V	74	THR	4.3
8	V	4	MET	4.2
2	D	12	THR	4.2
4	G	188	ASN	4.2
7	U	31(C)	TRP	4.1
7	U	31(B)	TYR	4.0
4	G	398	ASN	4.0
2	D	108	LEU	3.9
3	E	76	SER	3.9
8	V	86	TYR	3.8
5	H	182	VAL	3.8
2	D	42	GLY	3.8
2	D	84	SER	3.8
3	E	106(A)	LEU	3.8
8	V	54	LEU	3.8
8	V	46	LEU	3.8
8	V	78	LEU	3.8
7	U	84	SER	3.6
7	U	82(C)	LEU	3.6
8	V	11	LEU	3.6
8	V	58	VAL	3.6
2	D	82(B)	ASN	3.5
2	D	11	THR	3.5
3	E	79	ARG	3.5

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Mol	Chain	Res	Type	RSRZ
7	U	28	ASP	3.5
7	U	13	LYS	3.5
2	D	14	PRO	3.5
6	L	210	THR	3.5
7	U	8	GLY	3.4
6	L	7	TYR	3.4
3	E	98	PHE	3.4
5	H	204	LYS	3.4
2	D	43	ARG	3.4
2	D	46	GLU	3.4
7	U	82(A)	LYS	3.4
8	V	104	VAL	3.4
8	V	77	SER	3.2
4	G	396	ILE	3.2
8	V	37	GLN	3.2
7	U	90	TYR	3.1
5	H	73	THR	3.1
8	V	17	ASP	3.1
8	V	100	GLN	3.1
8	V	42	LYS	3.1
3	E	104	VAL	3.1
7	U	82	VAL	3.1
7	U	44	GLN	3.1
8	V	27	ARG	3.1
5	H	188	GLY	3.1
2	D	80	MET	3.0
5	H	206	ASP	2.9
2	D	38	ARG	2.9
5	H	193	ILE	2.9
7	U	16	ALA	2.9
8	V	1	ASP	2.9
7	U	110	VAL	2.9
8	V	62	PHE	2.8
5	H	136	LEU	2.8
5	H	157	LEU	2.8
2	D	88	GLY	2.8
8	V	64	GLY	2.8
2	D	67	VAL	2.8
2	D	89	THR	2.8
3	E	2	SER	2.8
8	V	14	SER	2.8
8	V	48	VAL	2.7

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Mol	Chain	Res	Type	RSRZ
7	U	45	LEU	2.7
8	V	71	PHE	2.6
5	H	205	VAL	2.6
6	L	70	ARG	2.6
2	D	44	GLY	2.6
2	D	83	THR	2.6
7	U	3	GLN	2.6
2	D	45	PRO	2.6
7	U	30	TYR	2.6
8	V	51	ALA	2.6
7	U	10	ALA	2.6
3	E	47	ILE	2.5
2	D	17	SER	2.5
7	U	61	TRP	2.5
4	G	505	VAL	2.5
8	V	72	SER	2.5
2	D	9	ALA	2.4
2	D	71	THR	2.4
3	E	69	TRP	2.4
4	G	397	SER	2.4
8	V	87	PHE	2.4
2	D	60	ALA	2.4
8	V	23	CYS	2.4
5	H	190	GLN	2.4
7	U	105	GLN	2.4
5	H	29	ILE	2.4
8	V	98	PHE	2.4
7	U	64	ASN	2.4
2	D	19	LYS	2.3
5	H	152	TRP	2.3
7	U	62	TYR	2.3
7	U	104	GLY	2.3
7	U	14	PRO	2.3
6	L	67(A)	ILE	2.3
5	H	196	VAL	2.3
8	V	84	ALA	2.2
5	H	123	ALA	2.2
7	U	7	SER	2.2
8	V	15	LEU	2.2
5	H	1	GLN	2.2
7	U	89	VAL	2.2
7	U	100(A)	SER	2.2

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Mol	Chain	Res	Type	RSRZ
7	U	111	VAL	2.2
2	D	57	LYS	2.2
8	V	39	LYS	2.2
7	U	27	ASP	2.2
7	U	31(D)	VAL	2.1
8	V	43	ALA	2.1
7	U	31	SER	2.1
2	D	5	VAL	2.1
4	G	412	ASP	2.1
7	U	9	ALA	2.1
5	H	202	ASN	2.1
3	E	12	SER	2.1
4	G	135	THR	2.1
5	H	189	THR	2.1
2	D	61	PRO	2.0
5	H	2	VAL	2.0
2	D	16	SER	2.0
7	U	15	GLY	2.0
3	E	15	LEU	2.0
2	D	107	THR	2.0
7	U	2	VAL	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, 95^{th} 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
11	MAN	F	5	11/12	0.65	0.36	79,108,132,134	0
12	MAN	I	4	11/12	0.67	0.35	82,116,125,126	0
11	BMA	F	3	11/12	0.71	0.30	86,116,128,129	0
12	BMA	I	3	11/12	0.77	0.25	106,110,120,120	0
10	NAG	C	2	14/15	0.78	0.60	89,114,120,120	0
11	MAN	F	4	11/12	0.78	0.20	74,91,108,113	0
10	NAG	J	1	14/15	0.79	0.28	50,85,107,121	0

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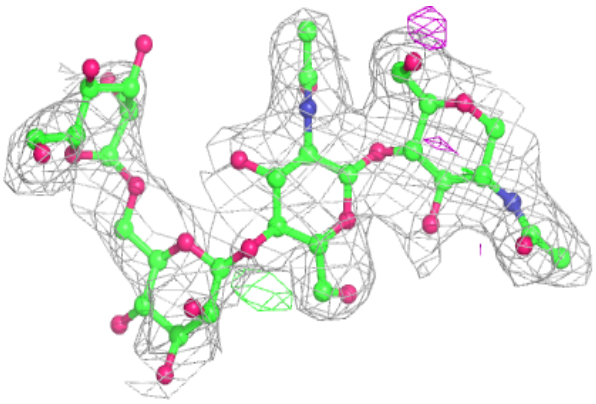
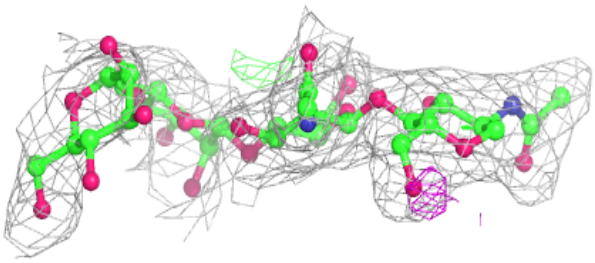
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
9	MAN	A	5	11/12	0.79	0.26	75,85,98,111	0
13	MAN	N	9	11/12	0.80	0.37	65,92,102,113	0
10	NAG	O	2	14/15	0.81	0.32	89,104,113,115	0
10	NAG	J	2	14/15	0.82	0.25	43,100,112,118	0
10	NAG	M	2	14/15	0.83	0.19	63,74,88,93	0
10	NAG	C	1	14/15	0.86	0.34	56,76,98,113	0
9	MAN	A	4	11/12	0.87	0.18	65,74,93,95	0
10	NAG	O	1	14/15	0.89	0.17	52,80,92,102	0
10	NAG	K	2	14/15	0.90	0.32	84,95,111,116	0
13	MAN	N	7	11/12	0.91	0.19	43,55,80,88	0
12	NAG	I	2	14/15	0.91	0.13	37,60,79,91	0
11	NAG	F	2	14/15	0.91	0.15	59,72,93,108	0
13	MAN	N	8	11/12	0.91	0.12	47,67,77,78	0
13	MAN	N	10	11/12	0.92	0.13	58,65,75,88	0
9	MAN	A	6	11/12	0.93	0.14	45,53,60,71	0
13	BMA	N	3	11/12	0.93	0.12	25,40,48,48	0
10	NAG	K	1	14/15	0.94	0.15	35,54,80,97	0
13	MAN	N	6	11/12	0.94	0.19	31,55,74,84	0
9	NAG	A	2	14/15	0.95	0.13	33,44,55,64	0
13	NAG	N	1	14/15	0.95	0.12	31,43,58,58	0
13	NAG	N	2	14/15	0.95	0.12	30,49,52,57	0
11	NAG	F	1	14/15	0.95	0.13	14,38,51,52	0
13	MAN	N	5	11/12	0.96	0.14	32,40,49,53	0
13	MAN	N	4	11/12	0.96	0.12	19,26,42,50	0
9	NAG	A	1	14/15	0.96	0.15	17,33,43,48	0
9	BMA	A	3	11/12	0.96	0.11	38,47,55,67	0
12	NAG	I	1	14/15	0.97	0.14	7,18,38,49	0
10	NAG	M	1	14/15	0.97	0.13	20,32,48,56	0

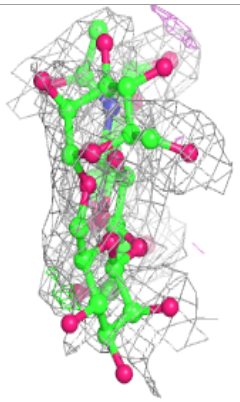
The following is a graphical depiction of the model fit to experimental electron density for oligosaccharide. Each fit is shown from different orientation to approximate a three-dimensional view.

Electron density around Chain I:

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

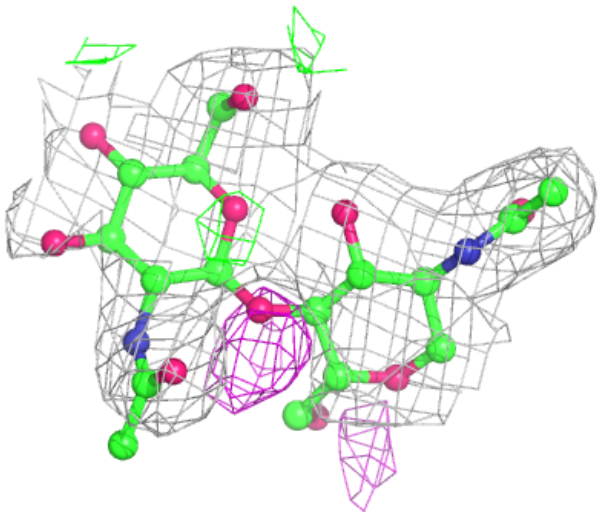


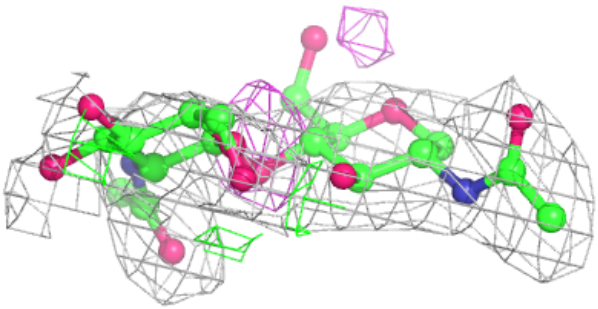


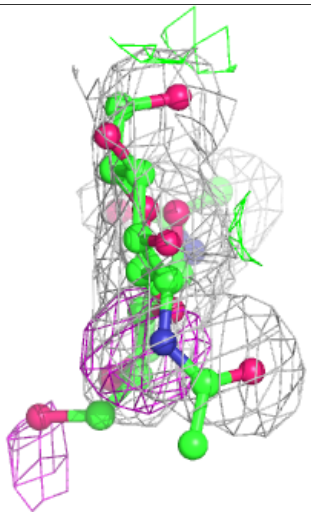


Electron density around Chain J:

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

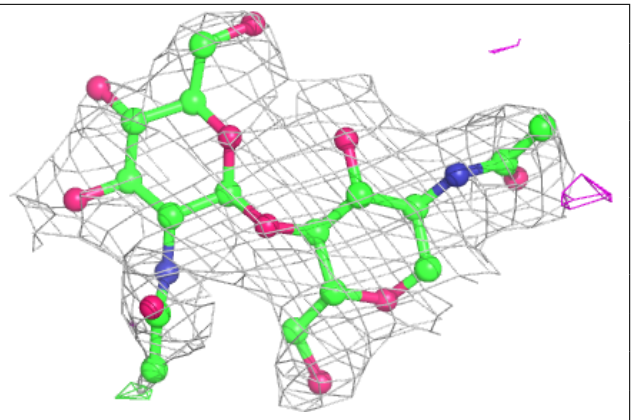


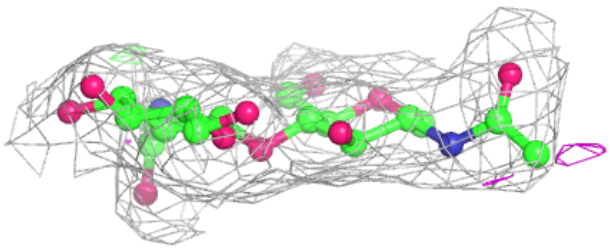


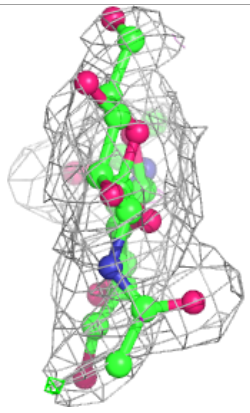


Electron density around Chain M:

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

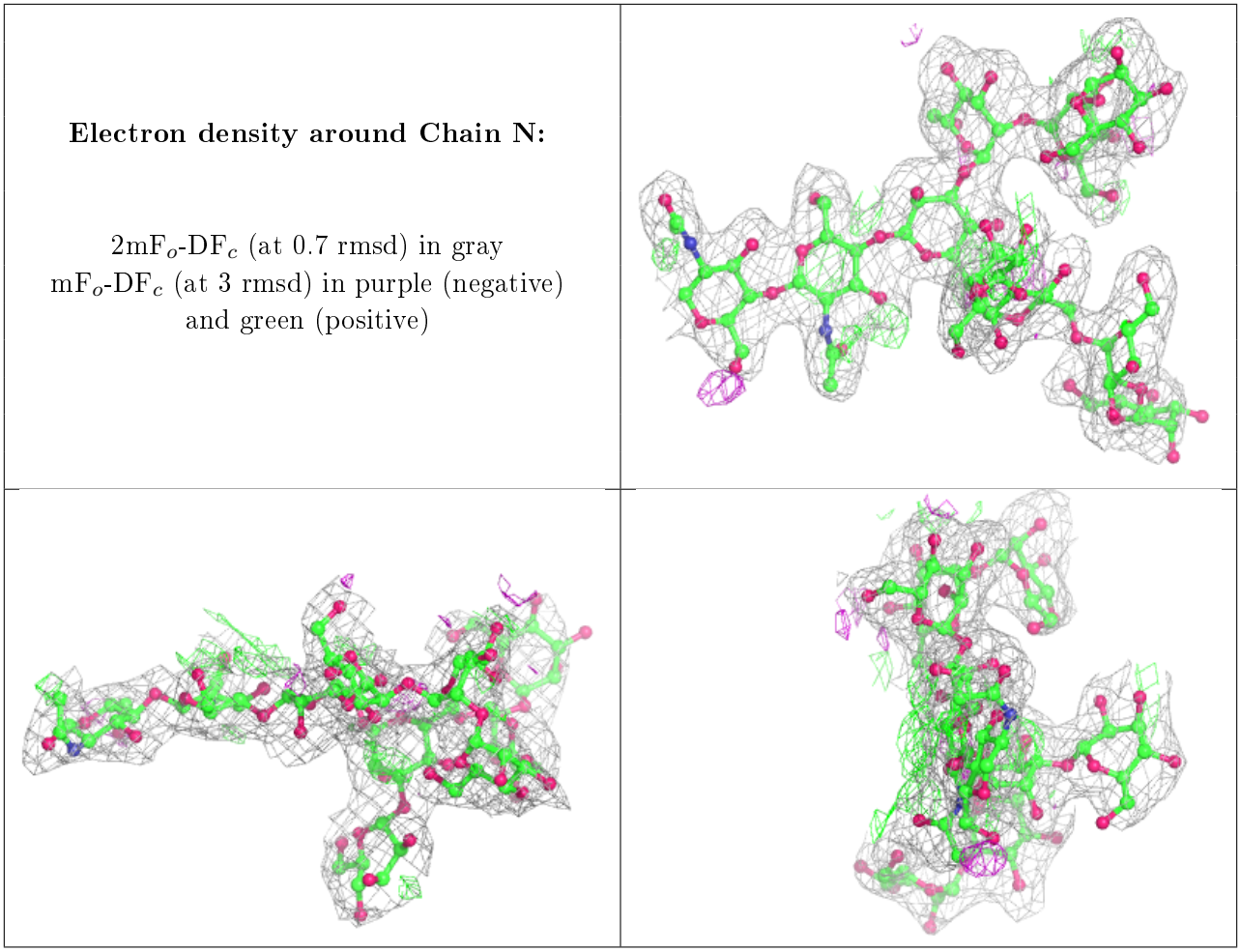






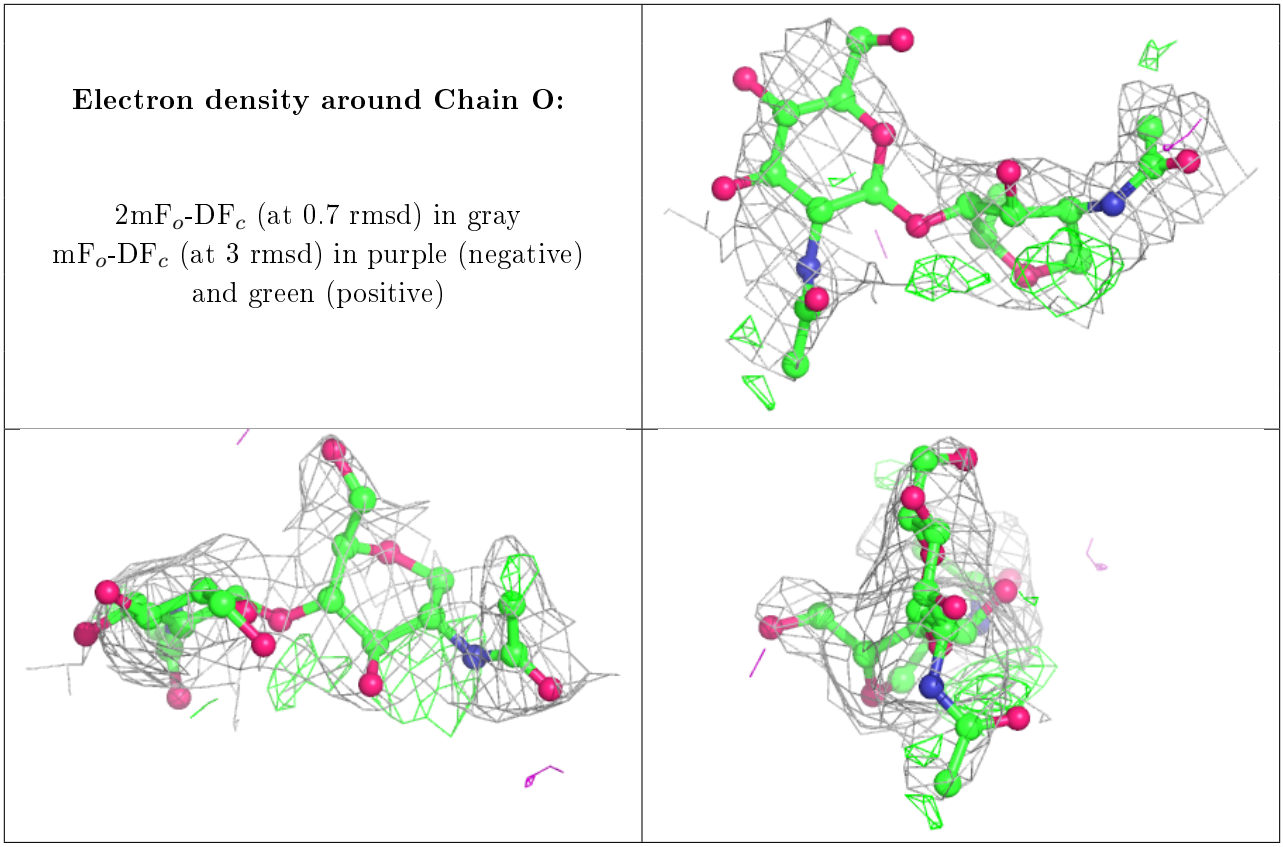
Electron density around Chain N:

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



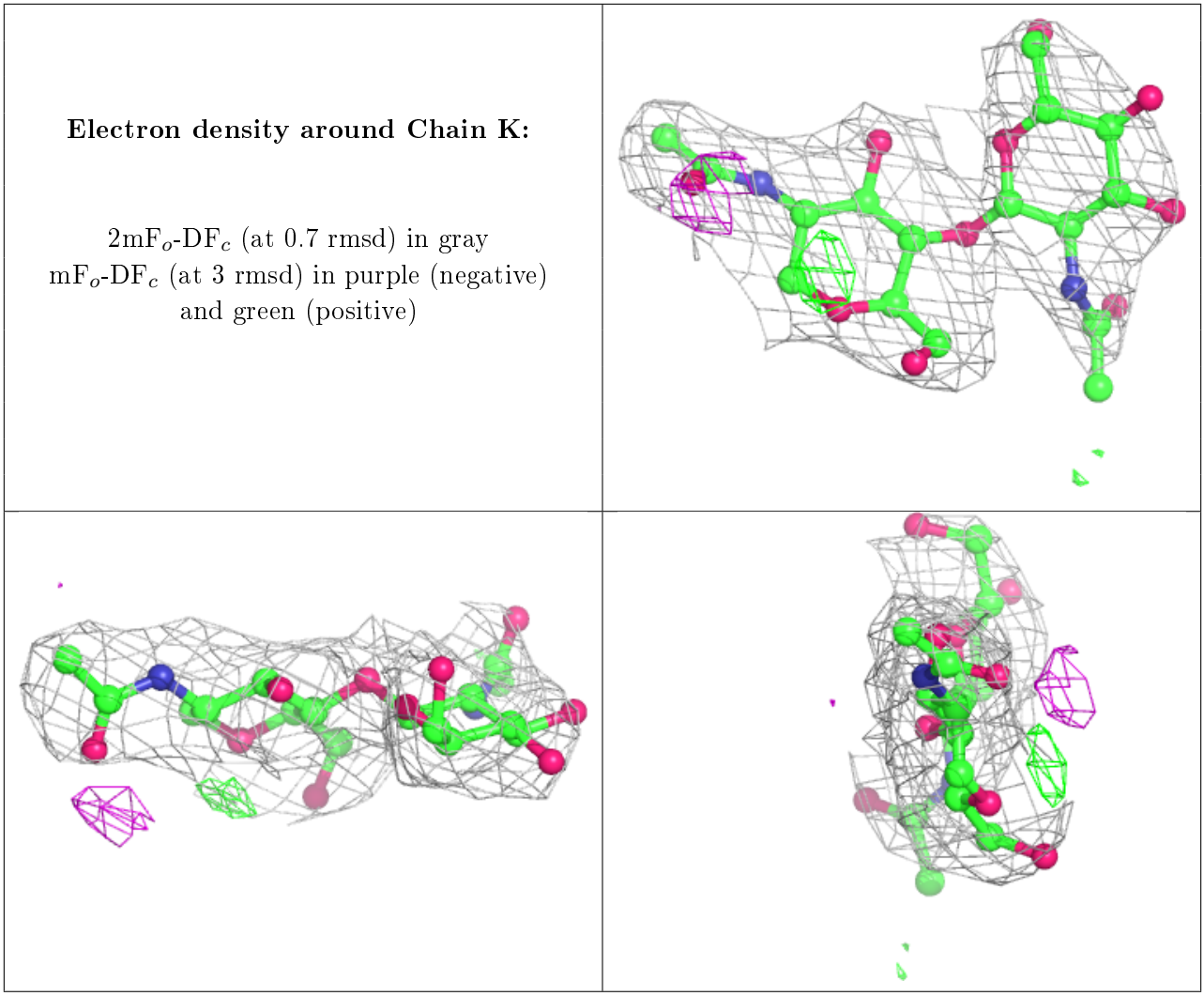
Electron density around Chain O:

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



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)



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, 95^{th} 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	G	616	14/15	0.81	0.22	66,88,101,107	0
14	NAG	B	702	14/15	0.84	0.31	78,92,101,102	0
14	NAG	G	615	14/15	0.88	0.15	37,60,79,82	0
14	NAG	B	701	14/15	0.89	0.44	72,99,112,124	0
14	NAG	G	637	14/15	0.92	0.14	36,47,59,63	0
14	NAG	G	638	14/15	0.93	0.14	36,42,51,66	0
14	NAG	G	614	14/15	0.93	0.13	21,55,84,89	0

6.5 Other polymers (i)

There are no such residues in this entry.