Full wwPDB X-ray Structure Validation Report (i)

Jun 22, 2024 – 02:56 PM EDT

PDB ID	:	4R0L
Title	:	Anti-canine CD28 antibody, 1C6, bound canine CD28
Authors	:	Gewe, M.M.; Strong, R.K.
Deposited on	:	2014-07-31
Resolution	:	3.30 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 types of validation reports are described at
http://www.wwpdb.org/validation/2017/FAQs#types.

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

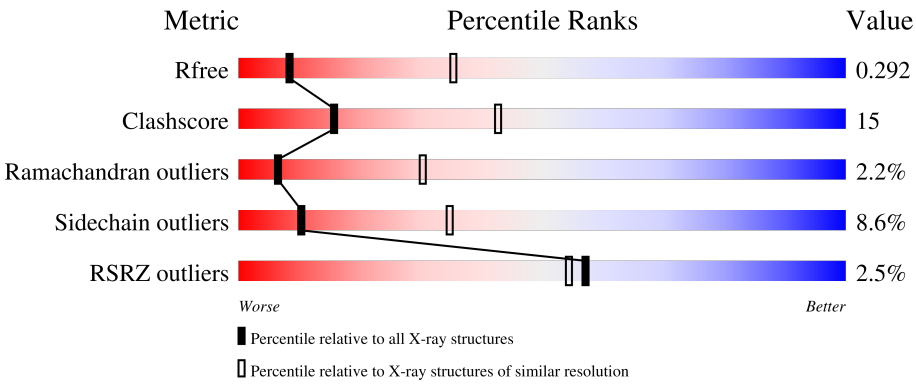
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.37.1
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.37.1

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 3.30 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	1149 (3.34-3.26)
Clashscore	141614	1205 (3.34-3.26)
Ramachandran outliers	138981	1183 (3.34-3.26)
Sidechain outliers	138945	1182 (3.34-3.26)
RSRZ outliers	127900	1115 (3.34-3.26)

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 <=5% The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	121	6% 70% 18% 5% 7%
1	H	121	2% 68% 21% • 8%
2	C	129	60% 24% 5% • 9%
2	D	129	% 63% 22% 5% • 9%
3	B	117	% 68% 29% •

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Mol	Chain	Length	Quality of chain
3	L	117	5% 73% 24%
4	E	3	100%
4	G	3	33% 67%
5	F	2	50% 50%

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
6	NAG	D	207	X	-	-	-

2 Entry composition (i)

There are 7 unique types of molecules in this entry. The entry contains 4887 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 Antibody 1C6 Heavy chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	113	746	471	126	145	4	0	0	0
1	H	111	738	476	121	137	4	0	0	0

- Molecule 2 is a protein called T-cell costimulatory molecule CD28.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	C	117	827	529	138	155	5	0	0	0
2	D	118	830	530	134	161	5	0	0	0

There are 22 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	0	SER	-	expression tag	UNP Q9N0N8
C	119	SER	-	expression tag	UNP Q9N0N8
C	120	ALA	-	expression tag	UNP Q9N0N8
C	121	TRP	-	expression tag	UNP Q9N0N8
C	122	SER	-	expression tag	UNP Q9N0N8
C	123	HIS	-	expression tag	UNP Q9N0N8
C	124	PRO	-	expression tag	UNP Q9N0N8
C	125	GLN	-	expression tag	UNP Q9N0N8
C	126	PHE	-	expression tag	UNP Q9N0N8
C	127	GLU	-	expression tag	UNP Q9N0N8
C	128	LYS	-	expression tag	UNP Q9N0N8
D	0	SER	-	expression tag	UNP Q9N0N8
D	119	SER	-	expression tag	UNP Q9N0N8
D	120	ALA	-	expression tag	UNP Q9N0N8
D	121	TRP	-	expression tag	UNP Q9N0N8
D	122	SER	-	expression tag	UNP Q9N0N8

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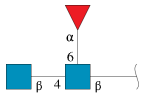
Chain	Residue	Modelled	Actual	Comment	Reference
D	123	HIS	-	expression tag	UNP Q9N0N8
D	124	PRO	-	expression tag	UNP Q9N0N8
D	125	GLN	-	expression tag	UNP Q9N0N8
D	126	PHE	-	expression tag	UNP Q9N0N8
D	127	GLU	-	expression tag	UNP Q9N0N8
D	128	LYS	-	expression tag	UNP Q9N0N8

- Molecule 3 is a protein called Antibody 1C6 Light chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
3	L	113	784	503	126	151	4	0	0	0
			Total	C	N	O	S			
3	B	114	786	503	130	149	4	0	0	0

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

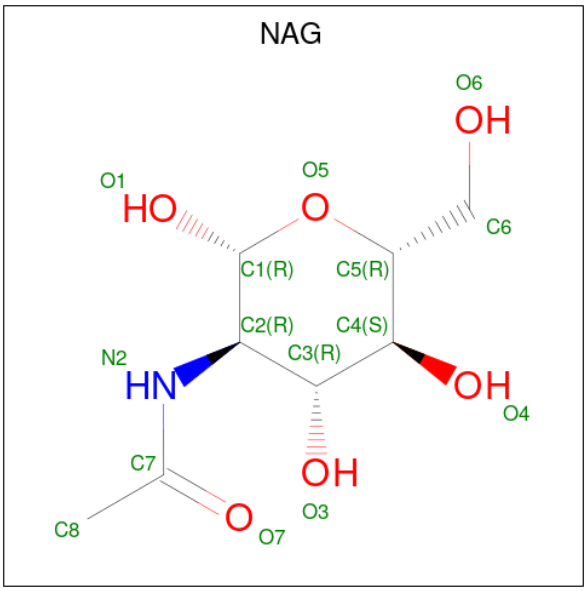


Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
			Total	C	N	O			
4	E	3	38	22	2	14	0	0	0
			Total	C	N	O			
4	G	3	38	22	2	14	0	0	0

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

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

- Molecule 6 is 2-acetamido-2-deoxy-beta-D-glucopyranose (three-letter code: NAG) (formula: $C_8H_{15}NO_6$).



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

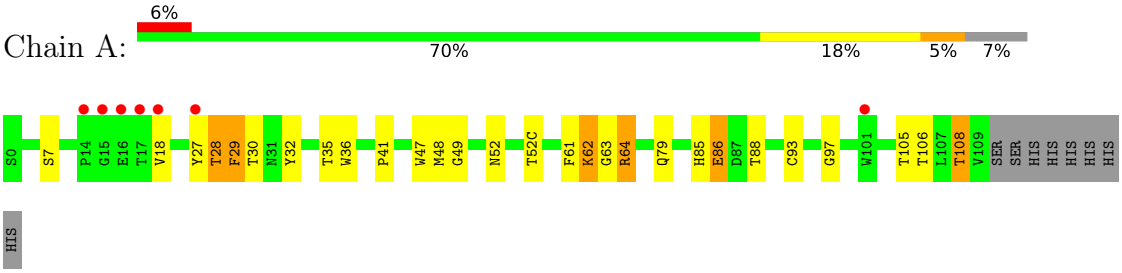
- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
7	C	1	1	1	0	0
			Total	O		
7	L	1	1	1	0	0

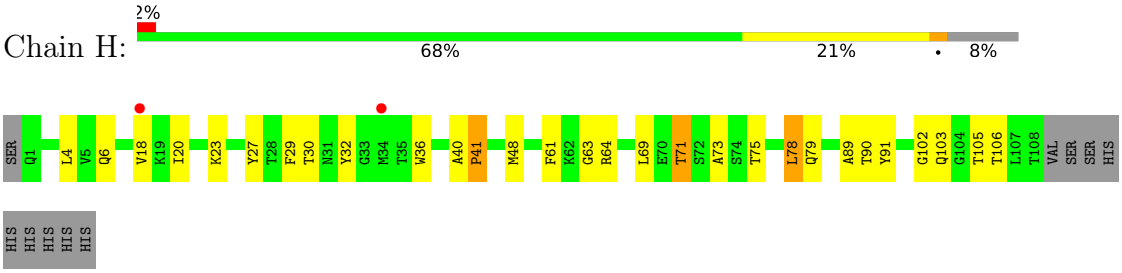
3 Residue-property plots (i)

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.

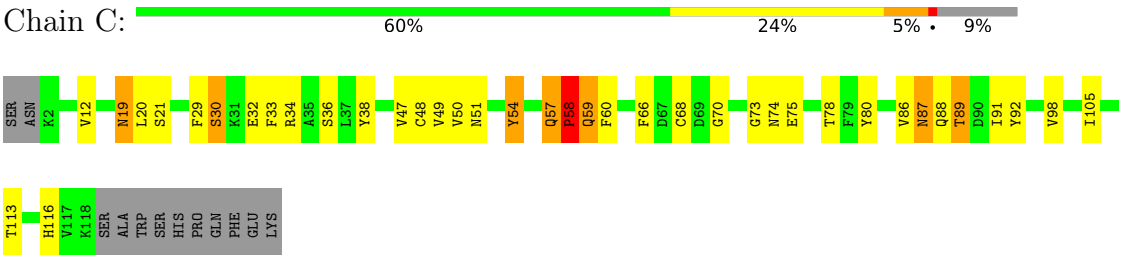
• Molecule 1: Antibody 1C6 Heavy chain



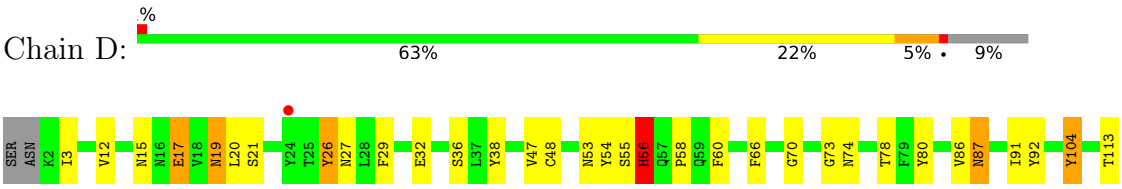
• Molecule 1: Antibody 1C6 Heavy chain



• Molecule 2: T-cell costimulatory molecule CD28

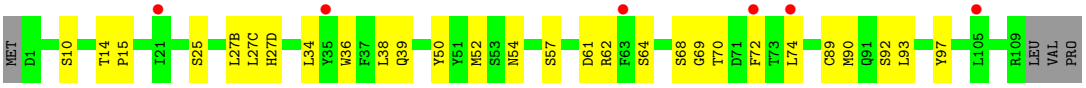


• Molecule 2: T-cell costimulatory molecule CD28



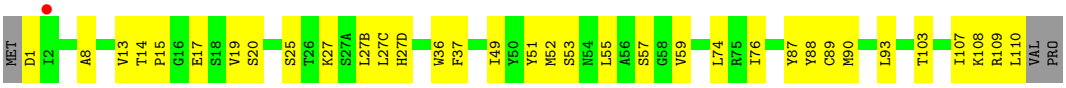
• Molecule 3: Antibody 1C6 Light chain

Chain L:



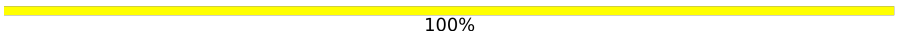
• Molecule 3: Antibody 1C6 Light chain

Chain B:



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

Chain E:



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

Chain G:

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

Chain F:

4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 31	Depositor
Cell constants a, b, c, α, β, γ	86.78Å 86.78Å 205.80Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	42.46 – 3.30 42.46 – 3.30	Depositor EDS
% Data completeness (in resolution range)	96.7 (42.46-3.30) 96.7 (42.46-3.30)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	0.08	Depositor
$< I/\sigma(I) >$ [1]	3.05 (at 3.32Å)	Xtriage
Refinement program	REFMAC 5.8.0049	Depositor
R, R_{free}	0.256 , 0.293 0.254 , 0.292	Depositor DCC
R_{free} test set	1292 reflections (5.11%)	wwPDB-VP
Wilson B-factor (Å^2)	105.8	Xtriage
Anisotropy	0.310	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.30 , 101.7	EDS
L-test for twinning[2]	$< \|L\| > = 0.52$, $< L^2 > = 0.36$	Xtriage
Estimated twinning fraction	0.003 for -h,-k,l 0.450 for h,-h-k,-l 0.007 for -k,-h,-l	Xtriage
F_o,F_c correlation	0.90	EDS
Total number of atoms	4887	wwPDB-VP
Average B, all atoms (Å^2)	95.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.44% 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 (i)

5.1 Standard geometry (i)

Bond lengths and bond angles in the following residue types are not validated in this section: FUC, NAG

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\#\|Z\|>5$	RMSZ	$\#\|Z\|>5$
1	A	0.54	0/767	0.62	0/1057
1	H	0.56	0/760	0.63	0/1046
2	C	0.59	0/849	0.69	0/1166
2	D	0.63	1/851 (0.1%)	0.69	0/1171
3	B	0.57	0/807	0.66	0/1109
3	L	0.56	0/805	0.66	0/1108
All	All	0.58	1/4839 (0.0%)	0.66	0/6657

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand.A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1
1	H	0	1
2	C	0	2
2	D	0	2
3	B	0	1
All	All	0	7

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	86	VAL	C-N	-5.80	1.20	1.34

There are no bond angle outliers.

There are no chirality outliers.

All (7) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	28	THR	Peptide
3	B	76	ILE	Peptide
2	C	54	TYR	Peptide
2	C	58	PRO	Peptide
2	D	118	LYS	Peptide
2	D	56	HIS	Peptide
1	H	69	LEU	Peptide

5.2 Too-close contacts (i)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	746	0	556	19	0
1	H	738	0	563	25	0
2	C	827	0	659	36	0
2	D	830	0	657	24	0
3	B	786	0	665	16	0
3	L	784	0	676	11	0
4	E	38	0	34	0	0
4	G	38	0	34	2	0
5	F	28	0	25	0	0
6	C	42	0	39	0	0
6	D	28	0	26	2	0
7	C	1	0	0	0	0
7	L	1	0	0	0	0
All	All	4887	0	3934	130	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 15.

All (130) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:52:ASN:ND2	2:C:59:GLN:OE1	1.97	0.95
1:H:73:ALA:O	1:H:75:THR:HG23	1.69	0.92
2:C:54:TYR:CD2	2:C:54:TYR:O	2.27	0.88
1:A:52:ASN:OD1	1:A:52(C):THR:CB	2.32	0.77

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:87:ASN:N	2:C:87:ASN:OD1	2.20	0.74
2:C:54:TYR:O	2:C:54:TYR:CG	2.39	0.73
2:D:19:ASN:N	2:D:19:ASN:OD1	2.23	0.72
3:L:50:TYR:O	3:L:54:ASN:HB2	1.91	0.71
1:H:78:LEU:HD12	1:H:79:GLN:N	2.07	0.70
2:C:29:PHE:HA	2:C:54:TYR:CD2	2.29	0.68
3:B:36:TRP:CZ3	3:B:89:CYS:HB3	2.31	0.66
2:C:12:VAL:HG23	2:C:116:HIS:HB3	1.78	0.65
2:D:21:SER:HA	2:D:78:THR:HA	1.80	0.64
1:A:27:TYR:CE2	1:A:29:PHE:HA	2.33	0.63
2:C:21:SER:HA	2:C:78:THR:HA	1.80	0.63
2:C:34:ARG:NH1	2:C:49:VAL:HG11	2.13	0.62
2:D:87:ASN:C	2:D:87:ASN:OD1	2.37	0.62
1:A:32:TYR:CE1	2:C:58:PRO:HB3	2.35	0.61
1:H:4:LEU:HD12	1:H:4:LEU:N	2.16	0.60
1:A:48:MET:HG2	1:A:61:PHE:CZ	2.37	0.59
1:A:48:MET:HG2	1:A:61:PHE:CE1	2.39	0.58
1:A:28:THR:O	1:A:29:PHE:CB	2.52	0.56
2:D:53:ASN:O	2:D:56:HIS:O	2.24	0.56
3:B:8:ALA:O	3:B:103:THR:HG23	2.07	0.55
2:C:89:THR:HB	2:C:116:HIS:HA	1.90	0.54
2:D:73:GLY:O	2:D:74:ASN:C	2.44	0.54
2:C:73:GLY:O	2:C:74:ASN:C	2.46	0.54
1:A:28:THR:O	1:A:29:PHE:HB3	2.09	0.53
2:D:70:GLY:HA2	2:D:80:TYR:HD1	1.74	0.53
2:C:51:ASN:OD1	2:C:51:ASN:C	2.47	0.53
3:L:25:SER:O	3:L:70:THR:OG1	2.27	0.53
3:B:49:ILE:HA	3:B:55:LEU:HA	1.91	0.53
2:D:3:ILE:HG21	2:D:26:TYR:CD1	2.44	0.53
2:D:118:LYS:O	2:D:119:SER:CB	2.57	0.52
1:H:78:LEU:HD12	1:H:78:LEU:C	2.29	0.52
2:C:51:ASN:OD1	2:C:51:ASN:O	2.27	0.52
1:A:18:VAL:O	1:A:79:GLN:HA	2.10	0.52
3:L:27(C):LEU:HG	3:L:27(D):HIS:N	2.23	0.52
1:H:48:MET:HG2	1:H:61:PHE:CZ	2.45	0.51
2:C:33:PHE:C	2:C:33:PHE:CD1	2.83	0.51
1:A:63:GLY:O	1:A:64:ARG:CB	2.58	0.51
3:L:61:ASP:OD1	3:L:62:ARG:N	2.44	0.50
3:B:36:TRP:CH2	3:B:89:CYS:HB3	2.46	0.50
1:H:6:GLN:OE1	1:H:102:GLY:HA3	2.11	0.50
2:C:91:ILE:HG13	2:C:113:THR:O	2.12	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:18:VAL:O	1:H:79:GLN:HA	2.11	0.49
1:H:27:TYR:CE1	1:H:29:PHE:HA	2.46	0.49
1:A:88:THR:HG23	1:A:88:THR:O	2.12	0.49
3:B:108:LYS:O	3:B:110:LEU:N	2.45	0.49
2:C:98:VAL:CG2	2:C:105:ILE:HG13	2.43	0.49
2:C:50:VAL:CG2	2:C:58:PRO:O	2.60	0.49
2:C:33:PHE:HE2	2:C:54:TYR:HB2	1.78	0.49
2:C:50:VAL:HG23	2:C:58:PRO:O	2.13	0.49
1:A:88:THR:OG1	1:A:108:THR:HA	2.12	0.49
2:C:33:PHE:HB3	2:C:98:VAL:HG12	1.95	0.49
2:C:47:VAL:HG13	2:C:66:PHE:CD2	2.48	0.49
1:H:63:GLY:O	1:H:64:ARG:CB	2.60	0.49
2:D:87:ASN:OD1	2:D:87:ASN:O	2.31	0.48
2:C:20:LEU:C	2:C:20:LEU:HD23	2.32	0.48
2:D:15:ASN:HB3	2:D:17:GLU:HB3	1.94	0.48
3:B:19:VAL:HG22	3:B:20:SER:N	2.29	0.48
1:A:47:TRP:CH2	1:A:49:GLY:HA2	2.48	0.48
1:A:30:THR:O	2:C:59:GLN:NE2	2.46	0.48
3:L:38:LEU:HD12	3:L:39:GLN:N	2.28	0.47
3:B:51:TYR:O	3:B:53:SER:N	2.46	0.47
2:C:88:GLN:O	2:C:92:TYR:OH	2.31	0.47
2:C:86:VAL:O	2:C:89:THR:HG23	2.15	0.47
2:D:27:ASN:C	2:D:29:PHE:H	2.18	0.47
6:D:207:NAG:O7	6:D:207:NAG:C3	2.62	0.47
4:G:3:FUC:H63	4:G:3:FUC:H2	1.96	0.47
2:C:50:VAL:HB	2:C:60:PHE:CD2	2.49	0.47
2:D:12:VAL:O	2:D:12:VAL:HG13	2.14	0.47
2:D:47:VAL:HG13	2:D:66:PHE:CG	2.49	0.47
2:D:58:PRO:HB3	1:H:32:TYR:CE1	2.49	0.47
2:D:12:VAL:HG23	2:D:116:HIS:HB3	1.97	0.46
1:H:18:VAL:O	1:H:18:VAL:HG13	2.15	0.46
1:H:6:GLN:H	1:H:103:GLN:NE2	2.14	0.46
1:H:36:TRP:O	1:H:48:MET:HB2	2.16	0.46
2:D:60:PHE:CE1	2:D:70:GLY:C	2.89	0.46
3:L:92:SER:HB2	3:L:97:TYR:CE2	2.50	0.46
2:C:34:ARG:CZ	2:C:49:VAL:HG11	2.45	0.46
3:B:15:PRO:HG3	3:B:107:ILE:HD11	1.98	0.46
2:D:104:TYR:CD1	2:D:104:TYR:C	2.89	0.46
2:C:70:GLY:HA2	2:C:80:TYR:HD1	1.80	0.45
1:A:61:PHE:O	1:A:62:LYS:C	2.55	0.45
1:H:6:GLN:O	1:H:103:GLN:NE2	2.48	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:B:14:THR:O	3:B:17:GLU:HB2	2.16	0.45
3:B:27(C):LEU:HG	3:B:27(D):HIS:N	2.31	0.45
2:C:36:SER:HA	2:C:48:CYS:O	2.16	0.45
1:H:23:LYS:CB	1:H:75:THR:HG22	2.46	0.45
1:H:90:THR:HA	1:H:106:THR:HA	1.99	0.45
2:D:29:PHE:HA	2:D:54:TYR:CD2	2.52	0.44
3:L:14:THR:HB	3:L:15:PRO:CD	2.47	0.44
3:L:34:LEU:HD22	3:L:72:PHE:CD2	2.52	0.44
1:A:7:SER:O	1:A:105:THR:HG23	2.17	0.44
3:B:27(B):LEU:HA	3:B:93:LEU:HD22	2.00	0.44
1:H:6:GLN:H	1:H:103:GLN:CD	2.19	0.44
2:C:66:PHE:CD1	2:C:66:PHE:N	2.85	0.44
1:A:35:THR:HG22	1:A:36:TRP:N	2.32	0.43
3:L:61:ASP:O	3:L:62:ARG:C	2.55	0.43
1:H:20:ILE:HG23	1:H:105:THR:HG21	2.00	0.43
3:B:25:SER:OG	3:B:27:LYS:O	2.30	0.43
1:A:85:HIS:O	1:A:86:GLU:C	2.57	0.43
1:H:30:THR:HG22	1:H:71:THR:CG2	2.49	0.43
1:H:89:ALA:HB3	1:H:91:TYR:CE1	2.54	0.43
1:H:4:LEU:N	1:H:4:LEU:CD1	2.81	0.43
1:H:6:GLN:H	1:H:103:GLN:HE22	1.66	0.43
2:D:92:TYR:CD1	2:D:115:ILE:CG1	3.02	0.43
3:B:13:VAL:O	3:B:108:LYS:HG2	2.18	0.43
2:C:19:ASN:N	2:C:19:ASN:OD1	2.52	0.42
2:C:20:LEU:C	2:C:20:LEU:CD2	2.87	0.42
2:D:19:ASN:HB3	2:D:80:TYR:CD2	2.54	0.42
6:D:207:NAG:O7	6:D:207:NAG:O3	2.33	0.42
2:C:29:PHE:HA	2:C:54:TYR:CE2	2.54	0.42
4:G:1:NAG:H62	4:G:3:FUC:H3	2.01	0.42
2:D:38:TYR:O	2:D:92:TYR:HA	2.19	0.42
1:H:30:THR:CG2	1:H:71:THR:CG2	2.98	0.42
1:A:97:GLY:HA3	2:C:57:GLN:HG2	2.01	0.42
3:L:36:TRP:CZ3	3:L:89:CYS:HB3	2.55	0.42
3:B:37:PHE:N	3:B:88:TYR:O	2.51	0.42
2:D:12:VAL:O	2:D:12:VAL:CG1	2.68	0.41
2:C:21:SER:OG	2:C:78:THR:HG23	2.20	0.41
2:D:56:HIS:N	2:D:56:HIS:CD2	2.88	0.41
3:L:27(B):LEU:HA	3:L:93:LEU:HD22	2.01	0.41
3:B:55:LEU:HD13	3:B:59:VAL:HB	2.03	0.41
2:C:38:TYR:O	2:C:92:TYR:HA	2.20	0.41
2:D:91:ILE:HA	2:D:113:THR:O	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:6:GLN:N	1:H:103:GLN:HE22	2.18	0.41
3:B:37:PHE:O	3:B:87:TYR:HA	2.22	0.40
1:H:40:ALA:HB1	1:H:41:PRO:CD	2.52	0.40

There are no symmetry-related clashes.

5.3 Torsion angles (i)

5.3.1 Protein backbone (i)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	111/121 (92%)	98 (88%)	8 (7%)	5 (4%)	2	15
1	H	109/121 (90%)	95 (87%)	13 (12%)	1 (1%)	17	48
2	C	115/129 (89%)	101 (88%)	11 (10%)	3 (3%)	5	27
2	D	116/129 (90%)	104 (90%)	11 (10%)	1 (1%)	17	48
3	B	112/117 (96%)	100 (89%)	10 (9%)	2 (2%)	8	35
3	L	111/117 (95%)	101 (91%)	7 (6%)	3 (3%)	5	26
All	All	674/734 (92%)	599 (89%)	60 (9%)	15 (2%)	6	30

All (15) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	62	LYS
2	C	59	GLN
1	A	29	PHE
3	L	68	SER
3	B	52	MET
3	B	109	ARG
1	A	86	GLU
2	C	30	SER
3	L	52	MET

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Mol	Chain	Res	Type
1	A	64	ARG
2	D	26	TYR
3	L	69	GLY
1	A	41	PRO
1	H	41	PRO
2	C	58	PRO

5.3.2 Protein sidechains (i)

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	51/102 (50%)	48 (94%)	3 (6%)	19	49
1	H	49/102 (48%)	47 (96%)	2 (4%)	30	61
2	C	71/119 (60%)	63 (89%)	8 (11%)	6	22
2	D	73/119 (61%)	62 (85%)	11 (15%)	3	13
3	B	67/101 (66%)	63 (94%)	4 (6%)	19	49
3	L	71/101 (70%)	66 (93%)	5 (7%)	15	43
All	All	382/644 (59%)	349 (91%)	33 (9%)	10	35

All (33) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	93	CYS
1	A	106	THR
1	A	108	THR
2	C	19	ASN
2	C	30	SER
2	C	32	GLU
2	C	57	GLN
2	C	68	CYS
2	C	75	GLU
2	C	87	ASN
2	C	89	THR
2	D	17	GLU

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Mol	Chain	Res	Type
2	D	19	ASN
2	D	20	LEU
2	D	32	GLU
2	D	36	SER
2	D	48	CYS
2	D	55	SER
2	D	56	HIS
2	D	87	ASN
2	D	104	TYR
2	D	116	HIS
1	H	71	THR
1	H	78	LEU
3	L	10	SER
3	L	57	SER
3	L	64	SER
3	L	74	LEU
3	L	90	MET
3	B	1	ASP
3	B	57	SER
3	B	74	LEU
3	B	90	MET

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
2	D	16	ASN
2	D	56	HIS

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)

8 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	$\#\|Z\| > 2$
4	NAG	E	1	4,2	14,14,15	0.35	0	17,19,21	1.01	2 (11%)
4	NAG	E	2	4	14,14,15	0.51	0	17,19,21	1.30	3 (17%)
4	FUC	E	3	4	10,10,11	0.50	0	14,14,16	0.90	1 (7%)
5	NAG	F	1	2,5	14,14,15	0.44	0	17,19,21	1.06	1 (5%)
5	NAG	F	2	5	14,14,15	0.31	0	17,19,21	0.78	0
4	NAG	G	1	4,2	14,14,15	0.39	0	17,19,21	1.72	1 (5%)
4	NAG	G	2	4	14,14,15	0.39	0	17,19,21	1.26	2 (11%)
4	FUC	G	3	4	10,10,11	0.58	0	14,14,16	2.05	5 (35%)

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	NAG	E	1	4,2	-	2/6/23/26	0/1/1/1
4	NAG	E	2	4	-	3/6/23/26	0/1/1/1
4	FUC	E	3	4	-	-	0/1/1/1
5	NAG	F	1	2,5	-	0/6/23/26	0/1/1/1
5	NAG	F	2	5	-	1/6/23/26	0/1/1/1
4	NAG	G	1	4,2	-	4/6/23/26	0/1/1/1
4	NAG	G	2	4	-	1/6/23/26	0/1/1/1
4	FUC	G	3	4	-	-	0/1/1/1

There are no bond length outliers.

All (15) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
4	G	1	NAG	C1-O5-C5	5.59	119.76	112.19
4	G	3	FUC	C1-C2-C3	5.11	115.95	109.67
4	E	2	NAG	O5-C1-C2	-3.13	106.34	111.29
4	G	2	NAG	C1-O5-C5	2.79	115.98	112.19

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
4	G	3	FUC	O5-C1-C2	2.67	114.89	110.77
4	G	2	NAG	C2-N2-C7	-2.66	119.11	122.90
4	E	2	NAG	C1-O5-C5	-2.62	108.64	112.19
4	E	3	FUC	O2-C2-C1	2.48	114.23	109.15
4	G	3	FUC	C1-O5-C5	2.40	118.23	112.78
4	E	2	NAG	C3-C4-C5	2.31	114.36	110.24
4	G	3	FUC	C3-C4-C5	2.31	113.37	109.77
4	E	1	NAG	O5-C5-C6	2.12	110.52	107.20
4	E	1	NAG	O5-C1-C2	-2.07	108.01	111.29
5	F	1	NAG	C4-C3-C2	2.05	114.02	111.02
4	G	3	FUC	O2-C2-C3	-2.02	106.09	110.14

There are no chirality outliers.

All (11) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	E	1	NAG	O5-C5-C6-O6
4	G	1	NAG	O5-C5-C6-O6
4	E	1	NAG	C4-C5-C6-O6
4	E	2	NAG	C4-C5-C6-O6
4	E	2	NAG	O5-C5-C6-O6
4	G	1	NAG	C8-C7-N2-C2
4	G	1	NAG	O7-C7-N2-C2
4	G	1	NAG	C4-C5-C6-O6
4	G	2	NAG	O5-C5-C6-O6
4	E	2	NAG	C3-C2-N2-C7
5	F	2	NAG	C3-C2-N2-C7

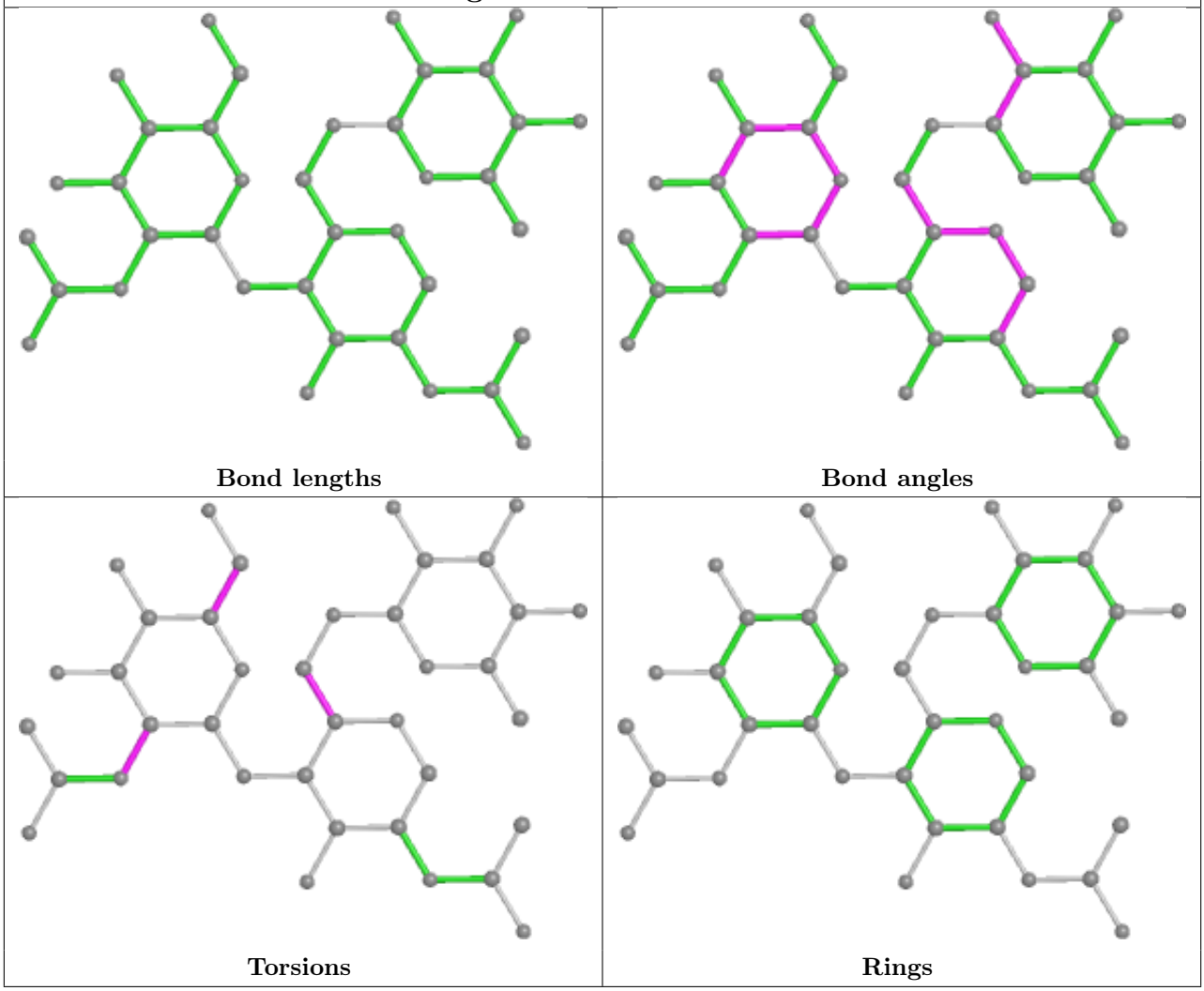
There are no ring outliers.

2 monomers are involved in 2 short contacts:

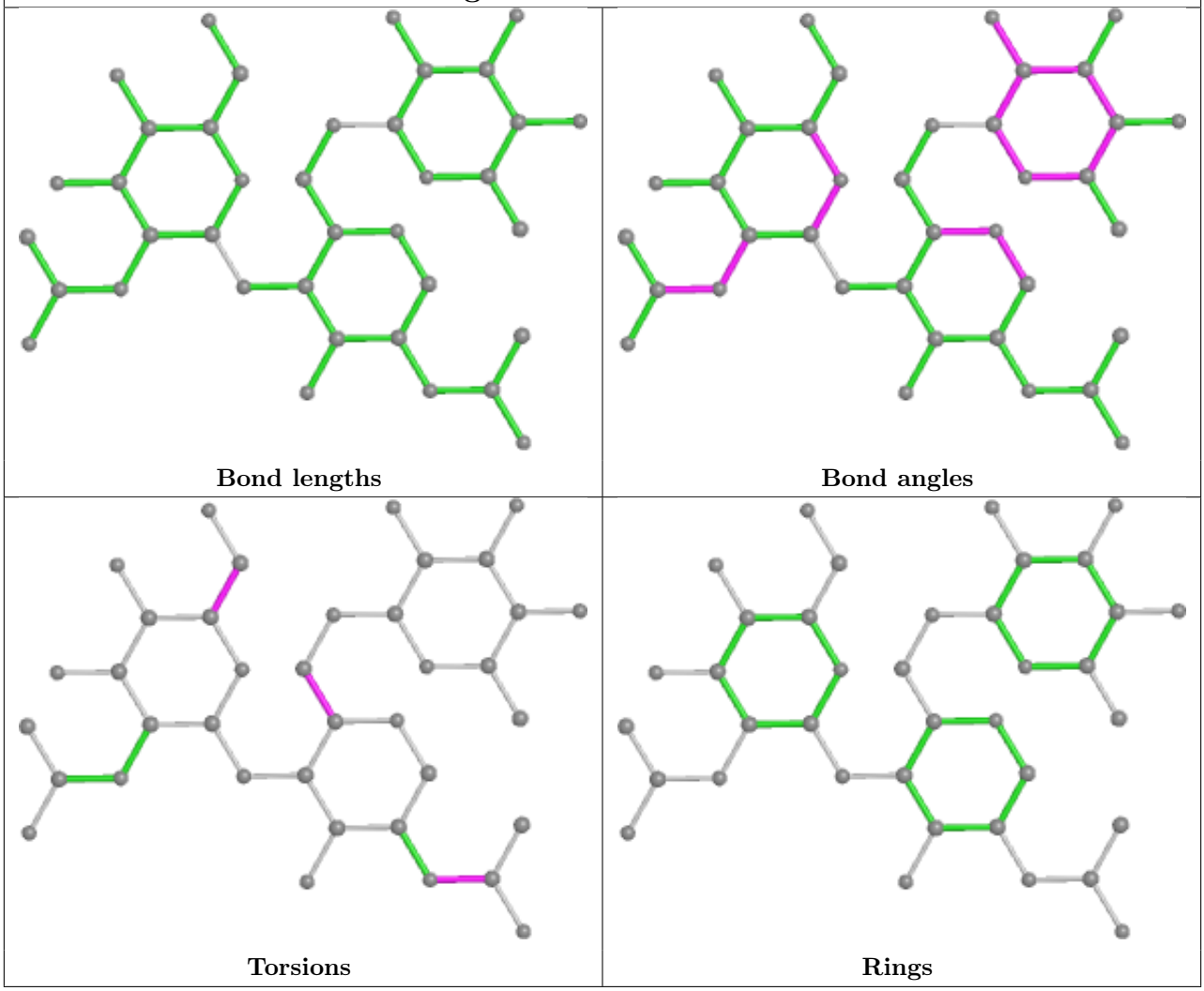
Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	G	1	NAG	1	0
4	G	3	FUC	2	0

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

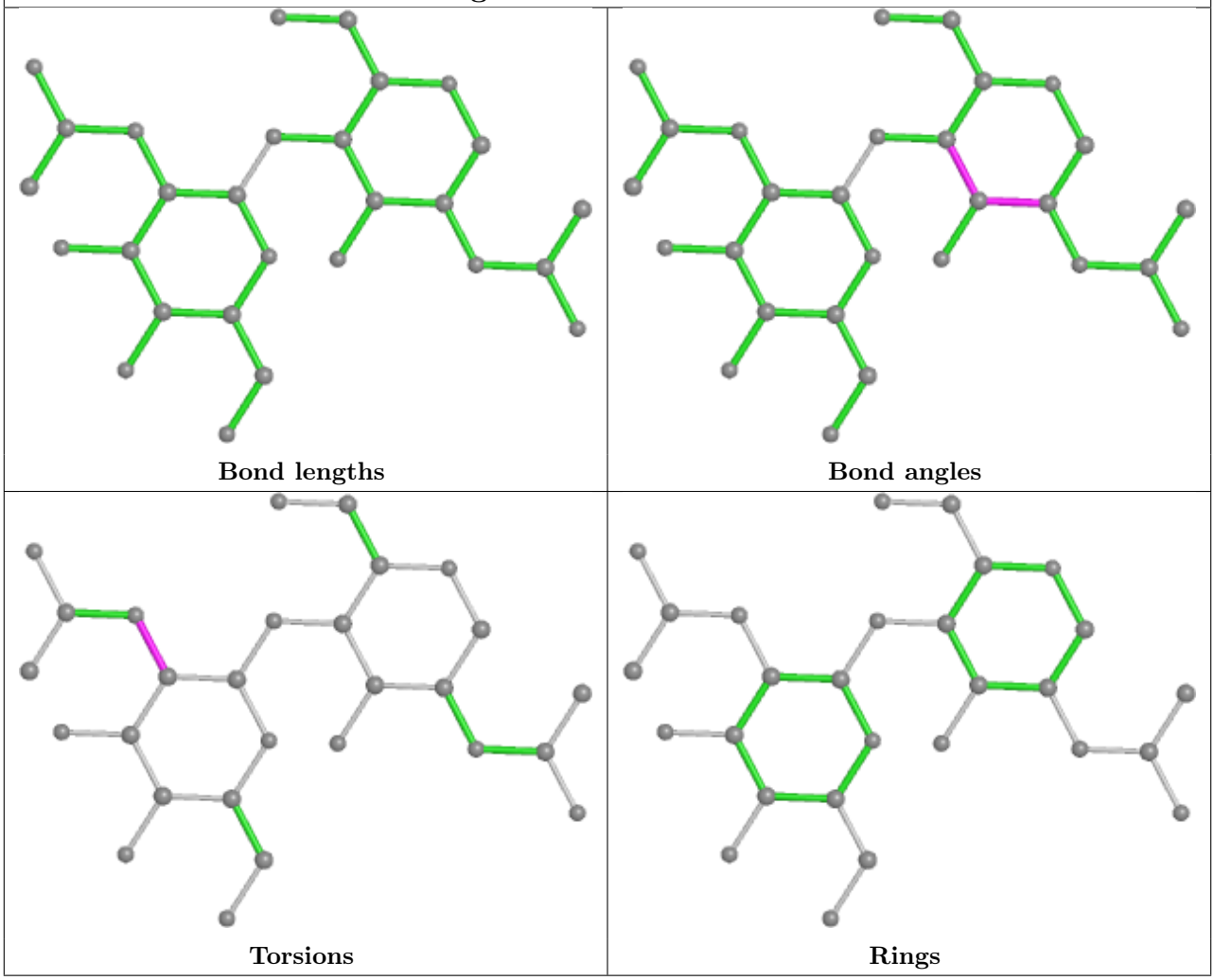
Oligosaccharide Chain E



Oligosaccharide Chain G



Oligosaccharide Chain F



5.6 Ligand geometry (i)

5 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	$\#\|Z\| > 2$
6	NAG	C	204	2	14,14,15	0.49	0	17,19,21	1.37	2 (11%)
6	NAG	C	206	2	14,14,15	0.45	0	17,19,21	1.40	3 (17%)
6	NAG	D	201	2	14,14,15	0.93	0	17,19,21	2.26	5 (29%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\| > 2$	Counts	RMSZ	$\#\|Z\| > 2$
6	NAG	D	207	2	14,14,15	0.49	0	17,19,21	1.61	4 (23%)
6	NAG	C	205	2	14,14,15	0.39	0	17,19,21	1.04	1 (5%)

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
6	NAG	C	204	2	-	2/6/23/26	0/1/1/1
6	NAG	C	206	2	-	2/6/23/26	0/1/1/1
6	NAG	D	201	2	-	4/6/23/26	0/1/1/1
6	NAG	D	207	2	1/1/5/7	1/6/23/26	0/1/1/1
6	NAG	C	205	2	-	2/6/23/26	0/1/1/1

There are no bond length outliers.

All (15) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
6	D	201	NAG	C1-O5-C5	5.53	119.69	112.19
6	D	207	NAG	C1-O5-C5	4.37	118.11	112.19
6	C	204	NAG	C1-O5-C5	3.99	117.60	112.19
6	D	201	NAG	C4-C3-C2	3.75	116.51	111.02
6	D	201	NAG	C2-N2-C7	3.50	127.89	122.90
6	C	206	NAG	C4-C3-C2	-3.27	106.22	111.02
6	D	201	NAG	O5-C1-C2	2.62	115.42	111.29
6	D	201	NAG	O5-C5-C6	2.54	111.18	107.20
6	C	206	NAG	O5-C1-C2	-2.52	107.31	111.29
6	C	204	NAG	O5-C1-C2	-2.42	107.47	111.29
6	D	207	NAG	C2-N2-C7	2.32	126.20	122.90
6	D	207	NAG	O5-C1-C2	2.28	114.89	111.29
6	C	205	NAG	O5-C5-C4	-2.14	105.62	110.83
6	C	206	NAG	C1-C2-N2	2.11	114.10	110.49
6	D	207	NAG	O5-C5-C6	2.10	110.50	107.20

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
6	D	207	NAG	C1

All (11) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
6	D	207	NAG	C3-C2-N2-C7
6	C	206	NAG	C4-C5-C6-O6
6	C	204	NAG	O5-C5-C6-O6
6	C	206	NAG	O5-C5-C6-O6
6	C	204	NAG	C4-C5-C6-O6
6	D	201	NAG	O5-C5-C6-O6
6	D	201	NAG	C8-C7-N2-C2
6	D	201	NAG	O7-C7-N2-C2
6	D	201	NAG	C4-C5-C6-O6
6	C	205	NAG	C4-C5-C6-O6
6	C	205	NAG	O5-C5-C6-O6

There are no ring outliers.

1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	D	207	NAG	2	0

5.7 Other polymers (i)

There are no such residues in this entry.

5.8 Polymer linkage issues (i)

There are no chain breaks in this entry.

6 Fit of model and data (i)

6.1 Protein, DNA and RNA chains (i)

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2			OWAB(Å^2)	Q<0.9
1	A	113/121 (93%)	0.17	7 (6%)	20	20	73, 99, 139, 157	0
1	H	111/121 (91%)	0.13	2 (1%)	68	67	73, 98, 133, 158	0
2	C	117/129 (90%)	0.00	0	100	100	62, 89, 110, 141	0
2	D	118/129 (91%)	0.18	1 (0%)	86	86	67, 89, 117, 133	0
3	B	114/117 (97%)	0.22	1 (0%)	84	84	72, 89, 118, 151	0
3	L	113/117 (96%)	0.23	6 (5%)	26	24	67, 91, 119, 144	0
All	All	686/734 (93%)	0.15	17 (2%)	57	54	62, 92, 124, 158	0

All (17) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	14	PRO	3.8
1	A	16	GLU	3.5
1	A	18	VAL	2.9
1	A	15	GLY	2.6
2	D	24	TYR	2.6
1	A	27	TYR	2.5
3	L	72	PHE	2.5
3	L	74	LEU	2.4
1	H	34	MET	2.4
3	L	35	TYR	2.3
3	L	21	ILE	2.2
1	A	17	THR	2.2
3	B	2	ILE	2.1
3	L	63	PHE	2.1
1	H	18	VAL	2.1
3	L	105	LEU	2.1
1	A	101	TRP	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)

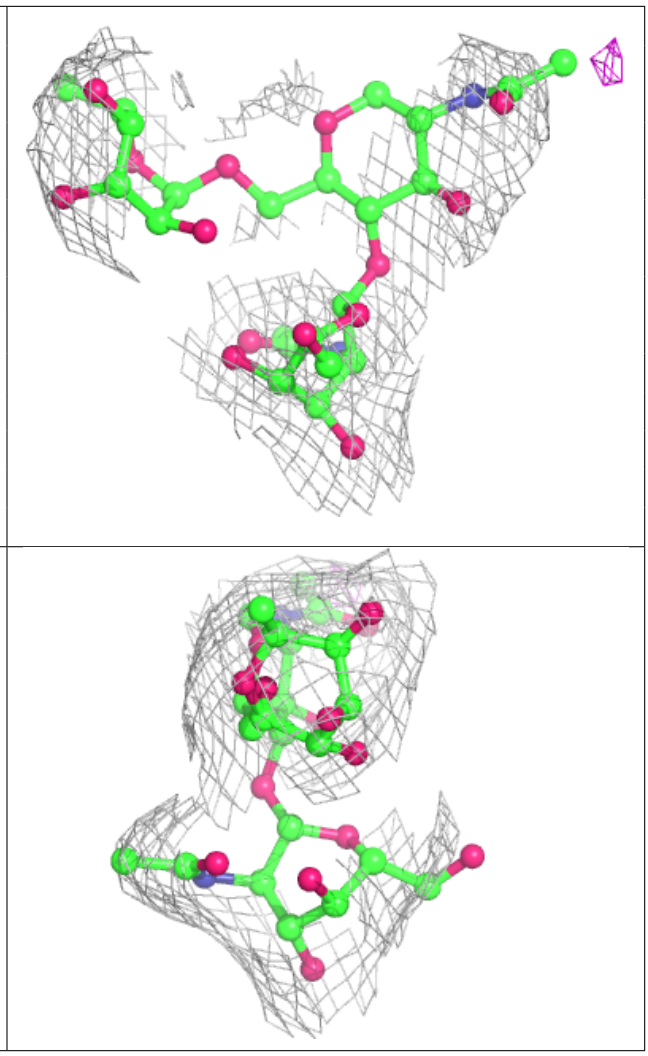
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
5	NAG	F	2	14/15	0.87	0.26	111,133,157,184	0
5	NAG	F	1	14/15	0.90	0.23	77,95,110,128	0
4	NAG	G	2	14/15	0.90	0.19	73,121,148,167	0
4	NAG	E	1	14/15	0.91	0.23	80,96,117,134	0
4	NAG	G	1	14/15	0.93	0.17	79,95,115,125	0
4	NAG	E	2	14/15	0.93	0.19	87,103,159,160	0
4	FUC	E	3	10/11	0.94	0.15	84,124,131,141	0
4	FUC	G	3	10/11	0.94	0.14	97,121,131,141	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.

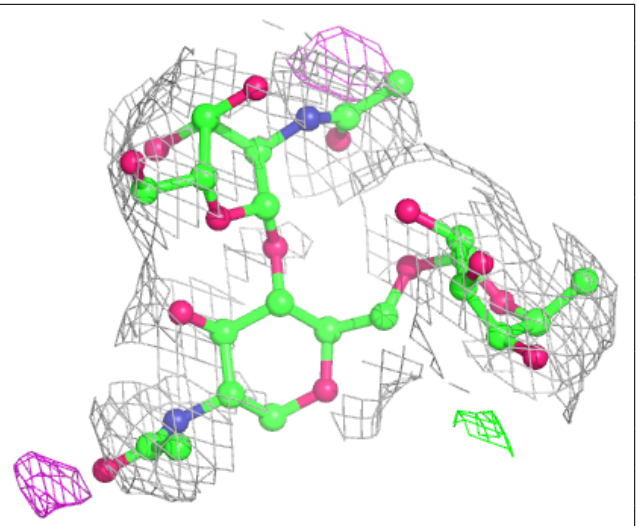
Electron density around Chain E:

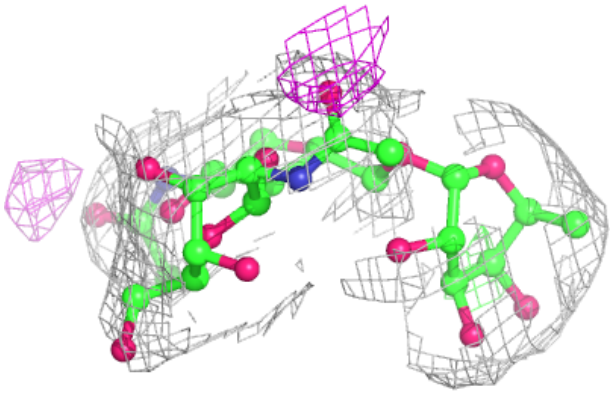
$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

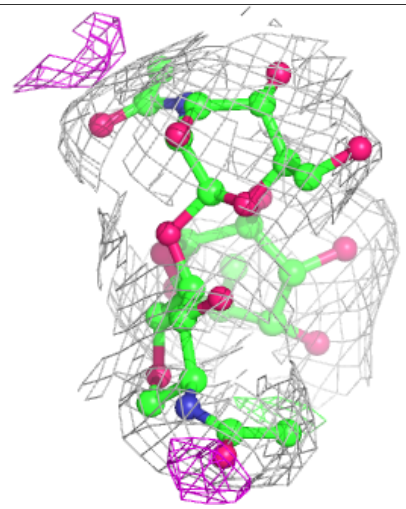


Electron density around Chain G:

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

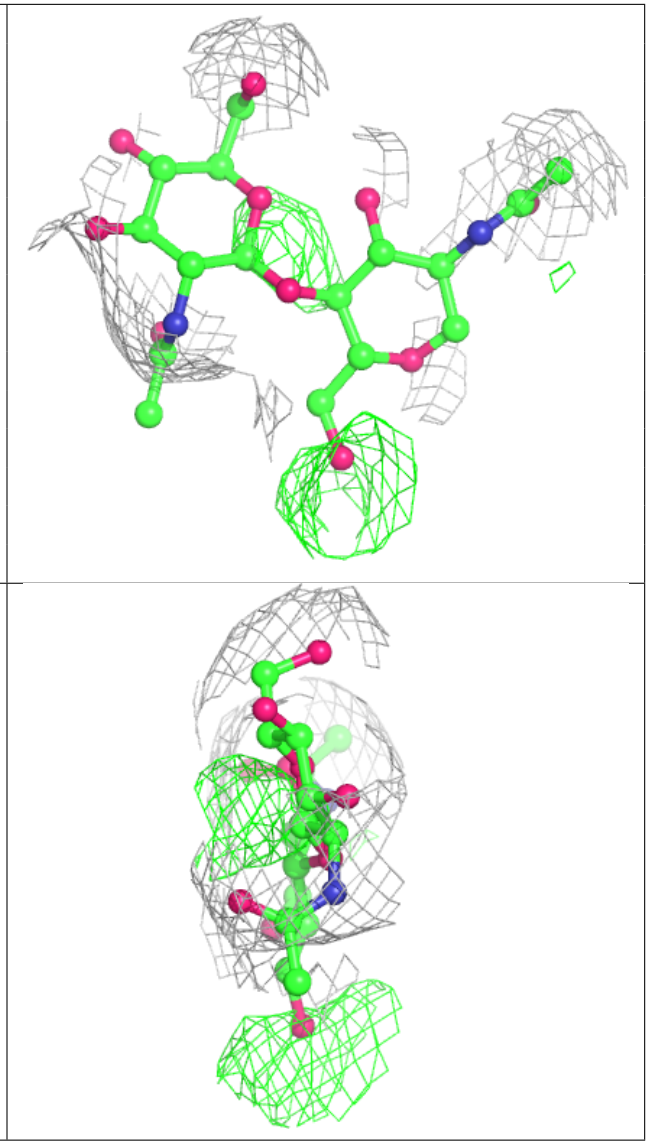






Electron density around Chain F:

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



6.4 Ligands (i)

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
6	NAG	C	205	14/15	0.57	0.26	77,103,122,138	0
6	NAG	D	201	14/15	0.69	0.28	65,90,187,206	0
6	NAG	C	204	14/15	0.81	0.27	71,86,110,140	0
6	NAG	C	206	14/15	0.86	0.17	100,137,162,187	0
6	NAG	D	207	14/15	0.87	0.18	85,105,130,131	0

6.5 Other polymers (i)

There are no such residues in this entry.