Full wwPDB X-ray Structure Validation Report ⓘ

Nov 20, 2023 – 01:05 AM JST

PDB ID	:	7C94
Title	:	Crystal structure of the anti-human podoplanin antibody Fab fragment complex with glycopeptide
Authors	:	Suzuki, K.; Nakamura, S.; Ogasawara, S.; Naruchi, K.; Shimabukuro, J.; Tukahara, N.; Kaneko, M.K.; Kato, Y.; Murata, T.
Deposited on	:	2020-06-04
Resolution	:	2.84 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:

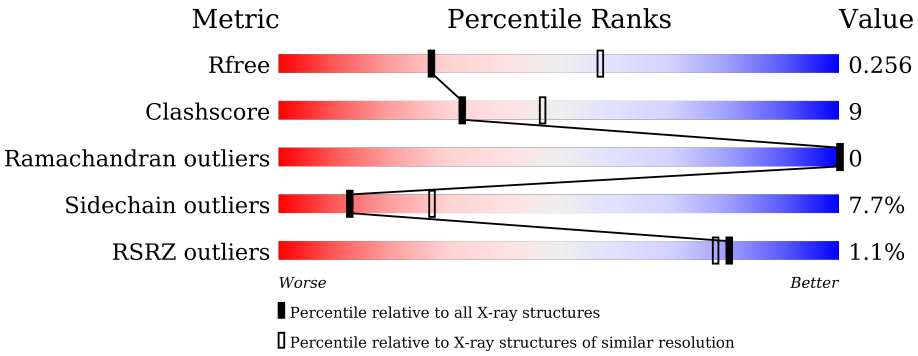
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.36
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36

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 2.84 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	1031 (2.86-2.82)
Clashscore	141614	1078 (2.86-2.82)
Ramachandran outliers	138981	1050 (2.86-2.82)
Sidechain outliers	138945	1051 (2.86-2.82)
RSRZ outliers	127900	1019 (2.86-2.82)

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	220	72% 27%
1	D	220	75% 23%
2	B	225	72% 25%
2	E	225	79% 19%
3	C	18	11% 28% 17% 56%
3	F	18	11% 17% 22% 6% 56%

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Mol	Chain	Length	Quality of chain
4	G	4	
4	H	4	

2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 7077 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 Light chain of Fab fragment.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	220	Total	C	N	O	S	0	0	0
			1696	1055	282	350	9			
1	D	219	Total	C	N	O	S	0	0	0
			1689	1051	281	349	8			

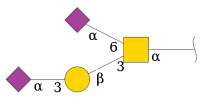
- Molecule 2 is a protein called Heavy chain of Fab fragment.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	223	Total	C	N	O	S	0	0	0
			1675	1061	281	324	9			
2	E	223	Total	C	N	O	S	0	0	0
			1680	1064	281	326	9			

- Molecule 3 is a protein called Peptide from Podoplanin.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
3	C	8	Total	C	N	O	0	0	0
			62	38	11	13			
3	F	8	Total	C	N	O	0	0	0
			62	38	11	13			

- Molecule 4 is an oligosaccharide called N-acetyl-alpha-neuraminic acid-(2-3)-beta-D-galacto pyranose-(1-3)-[N-acetyl-alpha-neuraminic acid-(2-6)]2-acetamido-2-deoxy-alpha-D-galacto pyranose.



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
4	G	4	Total	C	N	O	0	0	0
			65	36	3	26			

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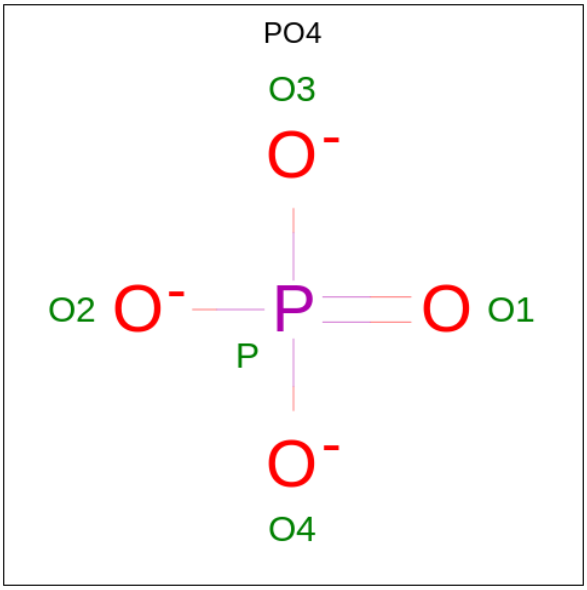
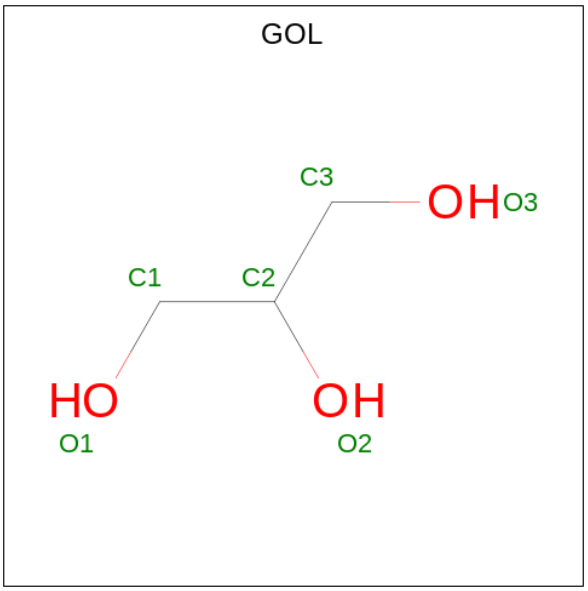
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf	Trace
4	H	4	Total C N O 65 36 3 26	0	0	0

- Molecule 5 is PHOSPHATE ION (three-letter code: PO4) (formula: O_4P).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	1	Total O P 5 4 1	0	0
5	D	1	Total O P 5 4 1	0	0
5	D	1	Total O P 5 4 1	0	0
5	E	1	Total O P 5 4 1	0	0

- Molecule 6 is GLYCEROL (three-letter code: GOL) (formula: $C_3H_8O_3$).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	A	1	Total C O 6 3 3	0	0
6	A	1	Total C O 6 3 3	0	0
6	B	1	Total C O 6 3 3	0	0
6	B	1	Total C O 6 3 3	0	0
6	D	1	Total C O 6 3 3	0	0
6	D	1	Total C O 6 3 3	0	0
6	D	1	Total C O 6 3 3	0	0

- Molecule 7 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
7	A	4	Total O 4 4	0	0
7	B	5	Total O 5 5	0	0
7	C	1	Total O 1 1	0	0
7	D	3	Total O 3 3	0	0
7	E	7	Total O 7 7	0	0

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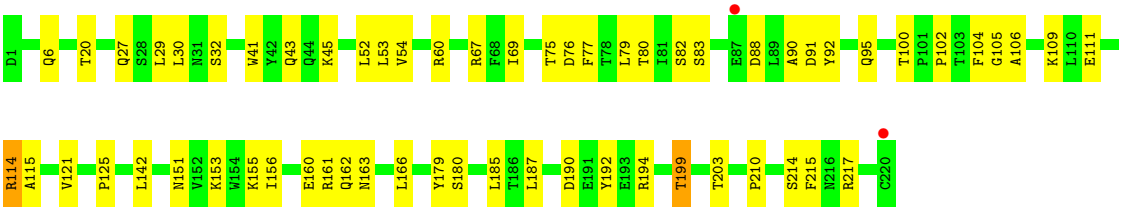
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
7	F	1	Total O 1 1	0	0

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.

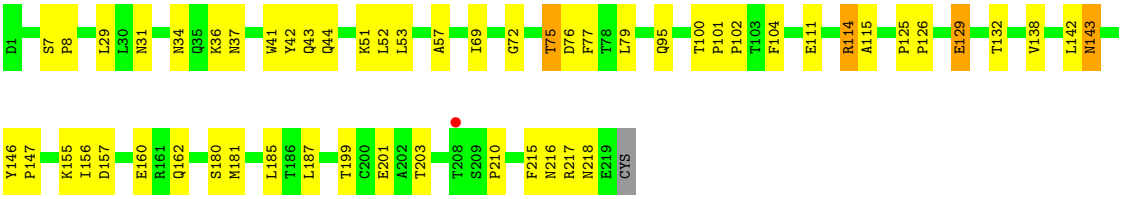
• Molecule 1: Light chain of Fab fragment

Chain A:



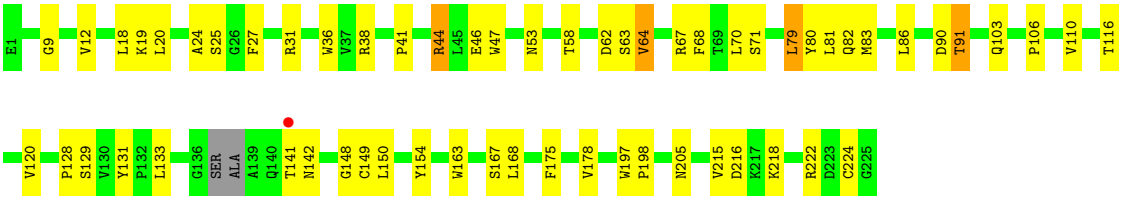
• Molecule 1: Light chain of Fab fragment

Chain D:



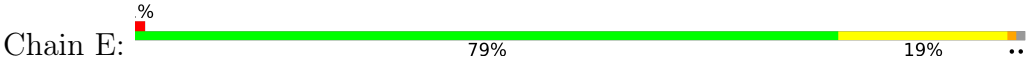
• Molecule 2: Heavy chain of Fab fragment

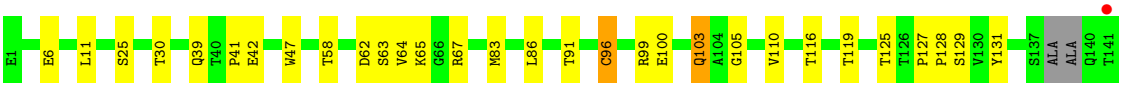
Chain B:

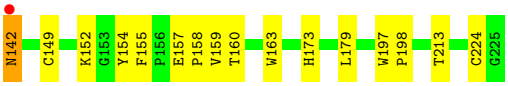


• Molecule 2: Heavy chain of Fab fragment

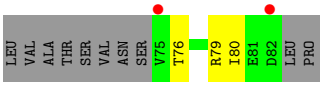
Chain E:



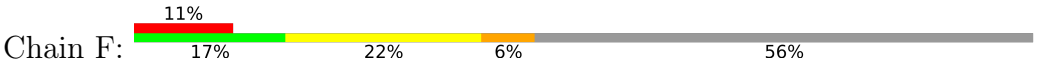


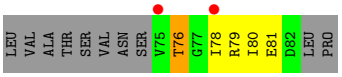


• Molecule 3: Peptide from Podoplanin



• Molecule 3: Peptide from Podoplanin





• Molecule 4: N-acetyl-alpha-neuraminic acid-(2-3)-beta-D-galactopyranose-(1-3)-[N-acetyl-alpha-neuraminic acid-(2-6)]2-acetamido-2-deoxy-alpha-D-galactopyranose

• Molecule 4: N-acetyl-alpha-neuraminic acid-(2-3)-beta-D-galactopyranose-(1-3)-[N-acetyl-alpha-neuraminic acid-(2-6)]2-acetamido-2-deoxy-alpha-D-galactopyranose

4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α, β, γ	43.10Å 42.96Å 221.08Å 90.00° 91.37° 90.00°	Depositor
Resolution (Å)	44.20 – 2.84 44.20 – 2.84	Depositor EDS
% Data completeness (in resolution range)	98.8 (44.20-2.84) 89.7 (44.20-2.84)	Depositor EDS
R_{merge}	0.18	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	0.90 (at 2.86Å)	Xtriage
Refinement program	PHENIX 1.14_3260	Depositor
R, R_{free}	0.220 , 0.255 0.220 , 0.256	Depositor DCC
R_{free} test set	1916 reflections (9.99%)	wwPDB-VP
Wilson B-factor (Å^2)	40.3	Xtriage
Anisotropy	0.457	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.37 , 23.1	EDS
L-test for twinning[2]	$< \|L\| > = 0.45$, $< L^2 > = 0.28$	Xtriage
Estimated twinning fraction	0.043 for k,h,-l 0.044 for -k,-h,-l 0.114 for h,-k,-l	Xtriage
F_o,F_c correlation	0.88	EDS
Total number of atoms	7077	wwPDB-VP
Average B, all atoms (Å^2)	46.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.25% 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: A2G, SIA, GOL, PO4, GAL

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	0.26	0/1733	0.44	0/2353
1	D	0.25	0/1726	0.44	0/2344
2	B	0.26	0/1718	0.46	0/2347
2	E	0.25	0/1723	0.46	0/2352
3	C	0.29	0/61	0.50	0/81
3	F	0.27	0/61	0.53	0/81
All	All	0.26	0/7022	0.45	0/9558

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1696	0	1614	33	0
1	D	1689	0	1608	33	0
2	B	1675	0	1630	34	0
2	E	1680	0	1640	22	0
3	C	62	0	62	3	0
3	F	62	0	62	3	0
4	G	65	0	53	2	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	H	65	0	53	8	0
5	A	5	0	0	0	0
5	D	10	0	0	0	0
5	E	5	0	0	0	0
6	A	12	0	16	0	0
6	B	12	0	16	1	0
6	D	18	0	24	1	0
7	A	4	0	0	0	0
7	B	5	0	0	0	0
7	C	1	0	0	1	0
7	D	3	0	0	0	0
7	E	7	0	0	0	0
7	F	1	0	0	1	0
All	All	7077	0	6778	120	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 (120) 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:D:37:ASN:HB2	1:D:57:ALA:HB2	1.71	0.73
1:A:156:ILE:HD11	1:A:185:LEU:HD21	1.71	0.72
2:E:142:ASN:OD1	2:E:142:ASN:N	2.18	0.71
2:E:83:MET:HB3	2:E:86:LEU:HD21	1.74	0.69
1:A:69:ILE:HB	1:A:80:THR:HB	1.75	0.68
1:D:52:LEU:HD22	2:E:110:VAL:HG22	1.74	0.68
2:E:91:THR:HG23	2:E:119:THR:HA	1.76	0.68
1:A:43:GLN:HB2	1:A:53:LEU:HD11	1.75	0.67
1:D:114:ARG:NH1	1:D:115:ALA:O	2.29	0.65
1:A:102:PRO:HG3	2:B:106:PRO:HB3	1.78	0.64
1:D:156:ILE:HD11	1:D:185:LEU:HD21	1.82	0.62
2:B:83:MET:HB3	2:B:86:LEU:HD21	1.81	0.62
2:B:131:TYR:HD2	2:B:150:LEU:HD23	1.65	0.62
2:E:128:PRO:HB3	2:E:154:TYR:HB3	1.83	0.61
2:E:63:SER:O	2:E:67:ARG:NH1	2.35	0.59
2:E:157:GLU:HB2	2:E:158:PRO:HA	1.84	0.59
3:F:79:ARG:NH2	7:F:301:HOH:O	2.34	0.59
2:E:99:ARG:NH2	2:E:105:GLY:O	2.33	0.59
1:D:31:ASN:HB3	1:D:34:ASN:HB2	1.85	0.59
1:A:67:ARG:NE	1:A:88:ASP:OD2	2.33	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:45:LYS:HD3	1:A:90:ALA:HB2	1.84	0.58
1:D:34:ASN:HB3	1:D:36:LYS:HG2	1.85	0.58
1:A:52:LEU:HD22	2:B:110:VAL:HG22	1.86	0.58
1:A:100:THR:OG1	4:G:4:SIA:O1B	2.15	0.57
3:C:79:ARG:NH1	7:C:301:HOH:O	2.37	0.57
2:B:41:PRO:HA	6:B:301:GOL:H2	1.86	0.56
1:D:129:GLU:O	1:D:132:THR:OG1	2.23	0.56
1:D:157:ASP:OD2	6:D:304:GOL:O2	2.24	0.56
4:H:2:GAL:H5	4:H:3:SIA:H113	1.89	0.55
2:B:38:ARG:NH2	2:B:46:GLU:OE2	2.41	0.54
1:A:125:PRO:HG2	2:B:222:ARG:CZ	2.38	0.54
2:B:128:PRO:HB3	2:B:154:TYR:HB3	1.90	0.53
4:H:4:SIA:O6	4:H:4:SIA:O8	2.25	0.53
1:A:106:ALA:HA	2:B:44:ARG:HE	1.74	0.53
1:D:101:PRO:HB3	2:E:47:TRP:CE3	2.43	0.52
2:B:36:TRP:NE1	2:B:79:LEU:HD13	2.24	0.52
2:B:53:ASN:N	2:B:53:ASN:OD1	2.43	0.52
1:D:95:GLN:HE21	1:D:102:PRO:HB3	1.75	0.52
2:B:91:THR:HG22	2:B:120:VAL:HB	1.92	0.52
1:D:143:ASN:HA	1:D:180:SER:HA	1.91	0.52
4:H:3:SIA:HN5	4:H:3:SIA:H92	1.73	0.52
2:B:67:ARG:NH2	2:B:90:ASP:OD2	2.44	0.51
1:D:43:GLN:HB2	1:D:53:LEU:HD11	1.92	0.50
1:D:142:LEU:HD13	1:D:181:MET:HG2	1.93	0.50
2:E:58:THR:O	4:H:3:SIA:O10	2.29	0.50
2:B:19:LYS:HE3	2:B:82:GLN:OE1	2.12	0.50
1:D:75:THR:O	1:D:75:THR:OG1	2.30	0.49
1:A:121:VAL:HG22	1:A:142:LEU:HD23	1.94	0.49
1:A:29:LEU:HD12	1:A:77:PHE:CE1	2.47	0.49
1:A:95:GLN:HB2	1:A:104:PHE:CD2	2.47	0.49
1:D:126:PRO:HD3	1:D:138:VAL:HG22	1.95	0.49
1:D:41:TRP:CE2	1:D:79:LEU:HB2	2.48	0.49
1:A:125:PRO:HB3	1:A:215:PHE:CE1	2.49	0.48
2:B:91:THR:HB	2:B:120:VAL:H	1.78	0.48
1:D:203:THR:O	1:D:203:THR:OG1	2.32	0.48
1:A:91:ASP:OD1	1:A:109:LYS:HG3	2.14	0.48
1:A:151:ASN:HB3	1:A:203:THR:OG1	2.13	0.48
4:G:2:GAL:H3	4:G:3:SIA:H32	1.44	0.47
2:E:100:GLU:HG3	2:E:110:VAL:HG23	1.97	0.47
1:A:114:ARG:HD3	1:A:115:ALA:O	2.15	0.47
1:A:190:ASP:O	1:A:194:ARG:HG3	2.14	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:63:SER:O	2:B:67:ARG:NH1	2.48	0.47
1:D:95:GLN:HB2	1:D:104:PHE:CD2	2.50	0.47
1:D:180:SER:HG	2:E:173:HIS:CE1	2.33	0.47
1:A:203:THR:HG22	1:A:210:PRO:HB3	1.97	0.46
1:D:217:ARG:NH1	1:D:218:ASN:HD21	2.14	0.46
1:A:155:LYS:HE2	1:A:160:GLU:HG2	1.97	0.45
1:D:72:GLY:HA3	1:D:77:PHE:HA	1.98	0.45
1:D:100:THR:N	4:H:4:SIA:O1A	2.50	0.45
2:E:62:ASP:HA	2:E:65:LYS:HE3	1.99	0.45
2:E:41:PRO:HD3	2:E:91:THR:O	2.16	0.45
4:H:3:SIA:C10	4:H:3:SIA:H7	2.46	0.45
1:A:192:TYR:CZ	1:A:217:ARG:HG3	2.52	0.45
1:A:6:GLN:HG3	1:A:105:GLY:HA3	1.99	0.44
2:B:133:LEU:HB2	2:B:148:GLY:C	2.38	0.44
1:A:88:ASP:O	1:A:92:TYR:OH	2.33	0.44
1:A:153:LYS:HE2	1:A:153:LYS:HB2	1.87	0.44
1:D:44:GLN:OE1	2:E:39:GLN:NE2	2.49	0.44
2:E:6:GLU:HG2	2:E:96:CYS:HB3	2.00	0.44
1:D:125:PRO:HB3	1:D:215:PHE:CE1	2.53	0.44
1:D:129:GLU:H	1:D:129:GLU:HG2	1.41	0.44
2:B:205:ASN:ND2	2:B:216:ASP:OD1	2.47	0.44
2:E:127:PRO:HB3	2:E:213:THR:HG21	2.00	0.44
4:H:3:SIA:H92	4:H:3:SIA:N5	2.33	0.44
1:A:166:LEU:HD22	2:B:178:VAL:HG11	1.99	0.43
2:E:149:CYS:HB2	2:E:163:TRP:CH2	2.54	0.43
1:D:43:GLN:O	1:D:51:LYS:N	2.47	0.43
2:B:64:VAL:HG13	2:B:68:PHE:HB2	2.00	0.43
1:D:29:LEU:HD12	1:D:77:PHE:CE1	2.54	0.43
1:D:114:ARG:HD3	1:D:115:ALA:O	2.17	0.43
1:A:54:VAL:HG22	1:A:60:ARG:HA	2.00	0.43
2:E:103:GLN:HB3	3:F:80:ILE:HB	2.01	0.43
3:F:76:THR:HG21	4:H:1:A2G:C7	2.48	0.43
1:A:41:TRP:CE2	1:A:79:LEU:HB2	2.55	0.42
1:A:180:SER:HB3	2:B:175:PHE:CE2	2.54	0.42
2:B:53:ASN:OD1	3:C:80:ILE:HA	2.19	0.42
2:B:103:GLN:HA	3:C:80:ILE:HB	2.00	0.42
1:D:42:TYR:HE1	1:D:95:GLN:HB3	1.84	0.42
1:A:111:GLU:OE1	1:A:179:TYR:OH	2.21	0.42
1:A:199:THR:HB	1:A:214:SER:OG	2.20	0.42
2:B:149:CYS:HB2	2:B:163:TRP:CH2	2.54	0.42
2:B:12:VAL:HG11	2:B:86:LEU:HD12	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:7:SER:HA	1:D:8:PRO:HA	1.82	0.42
1:A:102:PRO:HD2	2:B:47:TRP:CD2	2.56	0.41
2:B:9:GLY:HA2	2:B:18:LEU:HD21	2.01	0.41
1:A:153:LYS:HD2	1:A:160:GLU:OE2	2.20	0.41
2:B:163:TRP:HB2	2:B:168:LEU:HB2	2.02	0.41
2:E:129:SER:HB3	2:E:131:TYR:CZ	2.55	0.41
2:B:24:ALA:HB1	2:B:27:PHE:HE1	1.85	0.41
2:B:197:TRP:CG	2:B:198:PRO:HA	2.55	0.41
1:A:67:ARG:HB2	1:A:82:SER:O	2.20	0.41
2:B:20:LEU:HD12	2:B:81:LEU:HD23	2.01	0.41
2:B:197:TRP:CD1	2:B:198:PRO:HA	2.55	0.41
2:E:197:TRP:CG	2:E:198:PRO:HA	2.57	0.41
2:E:125:THR:HA	2:E:155:PHE:O	2.20	0.40
2:B:129:SER:HB3	2:B:131:TYR:CZ	2.56	0.40
1:D:146:TYR:CG	1:D:147:PRO:HA	2.56	0.40
1:D:155:LYS:NZ	1:D:201:GLU:OE2	2.53	0.40
1:D:203:THR:HG22	1:D:210:PRO:HB3	2.03	0.40
2:B:70:LEU:HD12	2:B:80:TYR:O	2.22	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	218/220 (99%)	213 (98%)	5 (2%)	0	100	100
1	D	217/220 (99%)	213 (98%)	4 (2%)	0	100	100
2	B	219/225 (97%)	217 (99%)	2 (1%)	0	100	100
2	E	219/225 (97%)	217 (99%)	2 (1%)	0	100	100
3	C	6/18 (33%)	6 (100%)	0	0	100	100
3	F	6/18 (33%)	6 (100%)	0	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
All	All	885/926 (96%)	872 (98%)	13 (2%)	0	100 100

There are no Ramachandran outliers to report.

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	194/196 (99%)	181 (93%)	13 (7%)	16 32
1	D	193/196 (98%)	181 (94%)	12 (6%)	18 35
2	B	184/188 (98%)	168 (91%)	16 (9%)	10 21
2	E	186/188 (99%)	172 (92%)	14 (8%)	13 28
3	C	7/16 (44%)	6 (86%)	1 (14%)	3 6
3	F	7/16 (44%)	4 (57%)	3 (43%)	0 0
All	All	771/800 (96%)	712 (92%)	59 (8%)	13 27

All (59) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	20	THR
1	A	27	GLN
1	A	30	LEU
1	A	32	SER
1	A	75	THR
1	A	76	ASP
1	A	83	SER
1	A	114	ARG
1	A	161	ARG
1	A	162	GLN
1	A	163	ASN
1	A	187	LEU
1	A	199	THR
2	B	25	SER

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Mol	Chain	Res	Type
2	B	31	ARG
2	B	44	ARG
2	B	58	THR
2	B	62	ASP
2	B	64	VAL
2	B	71	SER
2	B	79	LEU
2	B	91	THR
2	B	116	THR
2	B	141	THR
2	B	142	ASN
2	B	167	SER
2	B	215	VAL
2	B	218	LYS
2	B	224	CYS
3	C	76	THR
1	D	69	ILE
1	D	75	THR
1	D	76	ASP
1	D	111	GLU
1	D	114	ARG
1	D	129	GLU
1	D	143	ASN
1	D	160	GLU
1	D	162	GLN
1	D	187	LEU
1	D	199	THR
1	D	216	ASN
2	E	11	LEU
2	E	25	SER
2	E	30	THR
2	E	42	GLU
2	E	64	VAL
2	E	96	CYS
2	E	103	GLN
2	E	116	THR
2	E	142	ASN
2	E	152	LYS
2	E	159	VAL
2	E	160	THR
2	E	179	LEU
2	E	224	CYS

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Mol	Chain	Res	Type
3	F	76	THR
3	F	78	ILE
3	F	81	GLU

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

Mol	Chain	Res	Type
1	D	95	GLN
1	D	218	ASN

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	A2G	G	1	3,4	14,14,15	0.65	1 (7%)	17,19,21	1.57	1 (5%)
4	GAL	G	2	4	11,11,12	1.27	2 (18%)	15,15,17	1.38	2 (13%)
4	SIA	G	3	4	20,20,21	1.99	3 (15%)	24,28,31	2.44	6 (25%)
4	SIA	G	4	4	20,20,21	1.61	2 (10%)	24,28,31	1.91	4 (16%)
4	A2G	H	1	3,4	14,14,15	0.56	0	17,19,21	0.72	1 (5%)
4	GAL	H	2	4	11,11,12	0.44	0	15,15,17	1.01	0
4	SIA	H	3	4	20,20,21	1.95	3 (15%)	24,28,31	3.14	10 (41%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\|>2$	Counts	RMSZ	$\#\|Z\|>2$
4	SIA	H	4	4	20,20,21	1.68	2 (10%)	24,28,31	1.73	4 (16%)

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	A2G	G	1	3,4	-	2/6/23/26	0/1/1/1
4	GAL	G	2	4	-	1/2/19/22	0/1/1/1
4	SIA	G	3	4	-	10/18/34/38	0/1/1/1
4	SIA	G	4	4	-	4/18/34/38	0/1/1/1
4	A2G	H	1	3,4	-	2/6/23/26	0/1/1/1
4	GAL	H	2	4	-	0/2/19/22	0/1/1/1
4	SIA	H	3	4	-	11/18/34/38	0/1/1/1
4	SIA	H	4	4	-	7/18/34/38	0/1/1/1

All (13) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	G	3	SIA	C2-C1	6.68	1.58	1.52
4	H	4	SIA	C2-C1	6.38	1.58	1.52
4	G	4	SIA	C2-C1	5.74	1.57	1.52
4	H	3	SIA	C2-C1	5.51	1.57	1.52
4	H	3	SIA	O6-C2	4.45	1.49	1.43
4	G	3	SIA	O6-C2	4.23	1.49	1.43
4	H	3	SIA	C7-C6	3.25	1.57	1.53
4	G	4	SIA	O6-C2	2.40	1.47	1.43
4	G	3	SIA	C7-C6	2.40	1.56	1.53
4	G	2	GAL	C2-C3	2.36	1.56	1.52
4	G	2	GAL	O3-C3	2.34	1.48	1.43
4	H	4	SIA	O6-C2	2.27	1.46	1.43
4	G	1	A2G	C1-C2	2.03	1.55	1.52

All (28) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
4	H	3	SIA	C3-C4-C5	-10.40	98.89	111.46
4	G	4	SIA	C4-C3-C2	6.61	121.66	109.81
4	G	3	SIA	C6-O6-C2	6.50	125.23	111.34

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
4	H	3	SIA	C6-O6-C2	6.05	124.28	111.34
4	G	3	SIA	C3-C4-C5	-5.58	104.71	111.46
4	H	4	SIA	C4-C3-C2	5.43	119.54	109.81
4	G	3	SIA	C6-C5-N5	5.31	119.74	110.91
4	H	3	SIA	C4-C5-N5	4.90	120.08	110.38
4	G	1	A2G	C2-N2-C7	-4.64	116.29	122.90
4	H	3	SIA	O1A-C1-C2	-3.82	113.54	122.57
4	G	3	SIA	C4-C5-C6	-3.61	99.96	109.10
4	G	4	SIA	O1A-C1-C2	-3.49	114.32	122.57
4	H	3	SIA	C6-C5-N5	-3.18	105.64	110.91
4	H	4	SIA	O1A-C1-C2	-3.07	115.32	122.57
4	H	3	SIA	C4-C3-C2	2.97	115.12	109.81
4	G	4	SIA	C3-C4-C5	2.92	115.00	111.46
4	H	3	SIA	O6-C2-C3	2.91	114.47	110.46
4	H	3	SIA	C8-C7-C6	2.80	118.34	113.03
4	H	1	A2G	O5-C1-C2	-2.53	107.29	111.29
4	H	4	SIA	C6-O6-C2	2.44	116.57	111.34
4	H	3	SIA	O1B-C1-O1A	2.26	129.21	124.09
4	H	4	SIA	C3-C4-C5	2.22	114.15	111.46
4	G	3	SIA	O4-C4-C3	2.22	115.44	109.94
4	G	4	SIA	C6-O6-C2	2.18	116.00	111.34
4	G	2	GAL	C1-O5-C5	2.17	115.13	112.19
4	G	2	GAL	C1-C2-C3	2.15	112.31	109.67
4	G	3	SIA	O1A-C1-C2	-2.14	117.51	122.57
4	H	3	SIA	C9-C8-C7	2.02	116.80	112.41

There are no chirality outliers.

All (37) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	G	3	SIA	O7-C7-C8-C9
4	G	3	SIA	O7-C7-C8-O8
4	H	3	SIA	O1A-C1-C2-C3
4	H	3	SIA	O1B-C1-C2-C3
4	H	3	SIA	C5-C6-C7-C8
4	H	3	SIA	C5-C6-C7-O7
4	H	3	SIA	O6-C6-C7-C8
4	H	3	SIA	O6-C6-C7-O7
4	G	3	SIA	C6-C7-C8-C9
4	H	1	A2G	C4-C5-C6-O6
4	G	3	SIA	C11-C10-N5-C5
4	G	3	SIA	O10-C10-N5-C5

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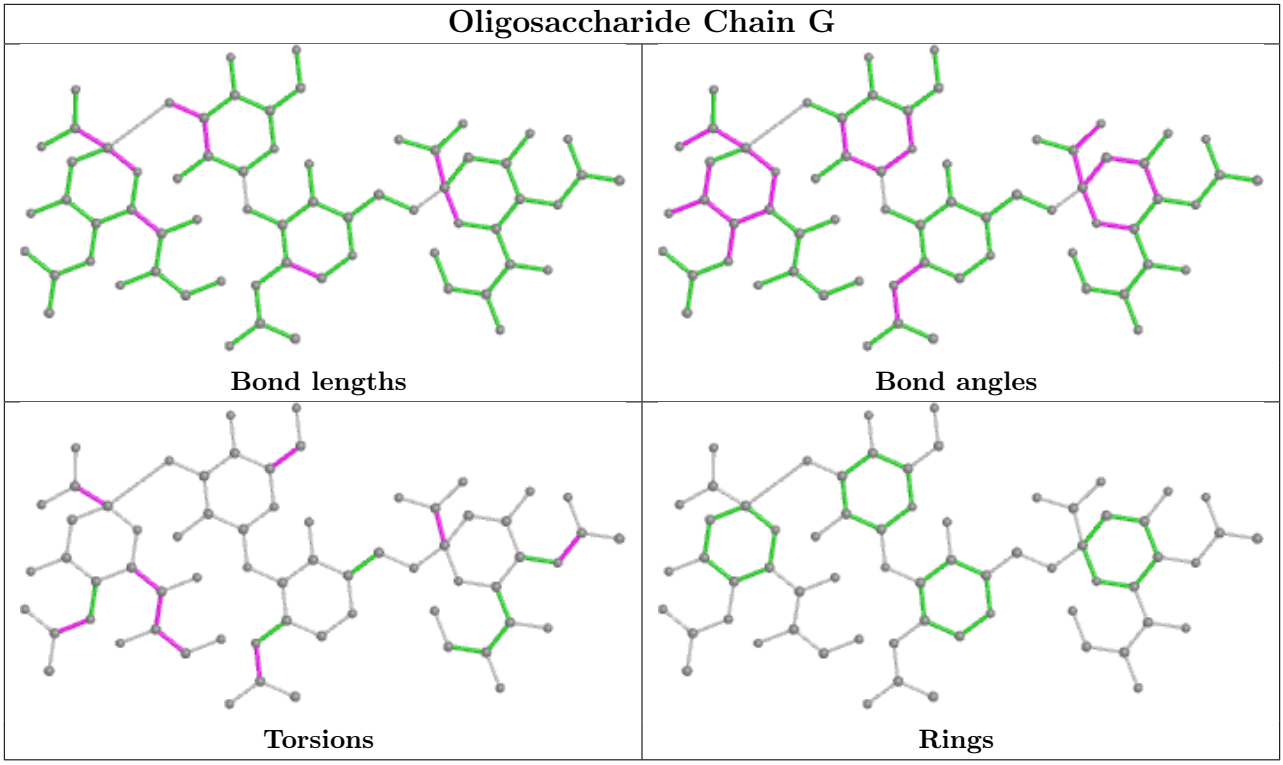
Mol	Chain	Res	Type	Atoms
4	G	4	SIA	C11-C10-N5-C5
4	G	4	SIA	O10-C10-N5-C5
4	H	3	SIA	C11-C10-N5-C5
4	H	3	SIA	O10-C10-N5-C5
4	H	4	SIA	C11-C10-N5-C5
4	H	4	SIA	O10-C10-N5-C5
4	H	3	SIA	O8-C8-C9-O9
4	H	1	A2G	O5-C5-C6-O6
4	G	3	SIA	C6-C7-C8-O8
4	H	4	SIA	O7-C7-C8-C9
4	H	4	SIA	C6-C7-C8-C9
4	H	3	SIA	O1A-C1-C2-O6
4	H	4	SIA	O1A-C1-C2-O6
4	H	4	SIA	O7-C7-C8-O8
4	G	3	SIA	C5-C6-C7-O7
4	H	3	SIA	C7-C8-C9-O9
4	G	1	A2G	C8-C7-N2-C2
4	G	3	SIA	C7-C8-C9-O9
4	G	3	SIA	O1A-C1-C2-O6
4	G	4	SIA	O1A-C1-C2-O6
4	G	3	SIA	O8-C8-C9-O9
4	G	4	SIA	O1B-C1-C2-C3
4	G	1	A2G	O7-C7-N2-C2
4	H	4	SIA	C6-C7-C8-O8
4	G	2	GAL	C4-C5-C6-O6

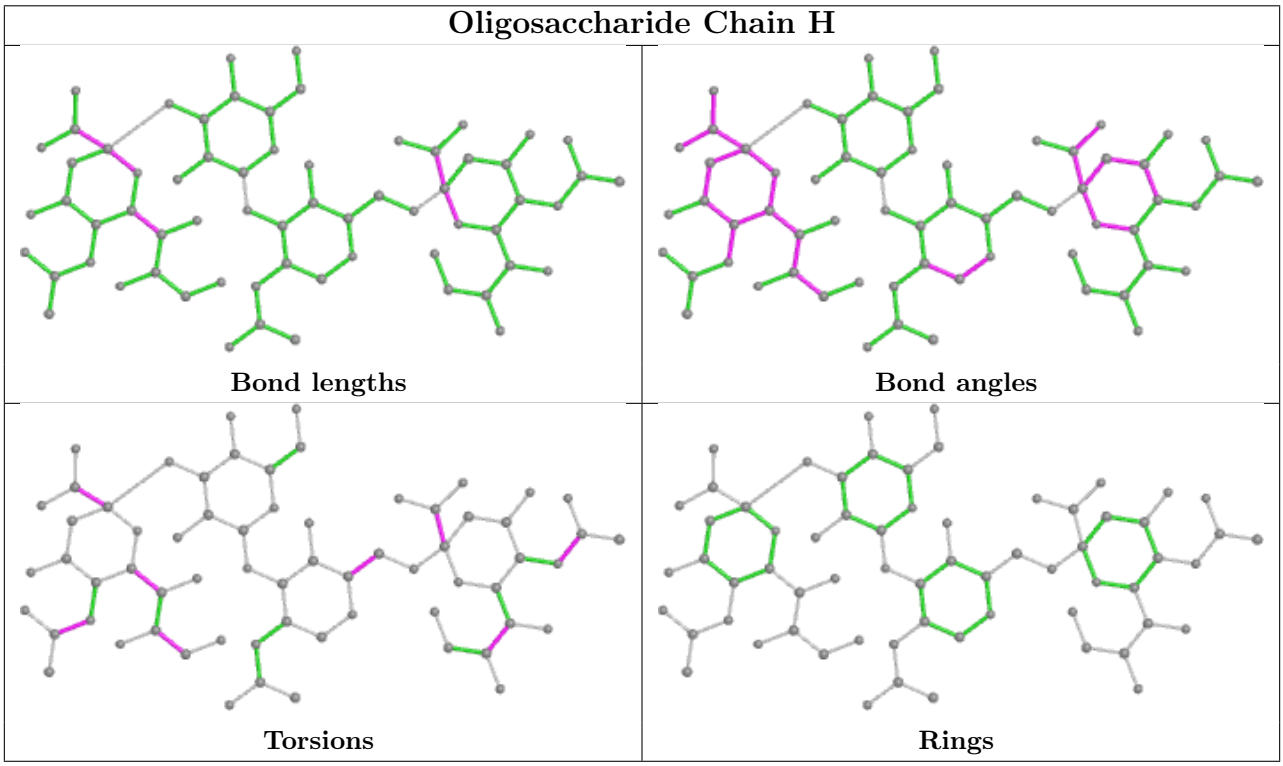
There are no ring outliers.

7 monomers are involved in 10 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	G	2	GAL	1	0
4	G	3	SIA	1	0
4	H	1	A2G	1	0
4	H	3	SIA	5	0
4	H	4	SIA	2	0
4	G	4	SIA	1	0
4	H	2	GAL	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 (i)

11 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	$\#\|Z\| > 2$
5	PO4	A	301	-	4,4,4	0.92	0	6,6,6	0.43	0
6	GOL	D	304	-	5,5,5	0.93	0	5,5,5	0.97	0
6	GOL	A	302	-	5,5,5	0.89	0	5,5,5	1.01	0
6	GOL	B	301	-	5,5,5	0.88	0	5,5,5	1.05	0
6	GOL	D	303	-	5,5,5	0.91	0	5,5,5	0.97	0
5	PO4	D	301	-	4,4,4	0.92	0	6,6,6	0.45	0
6	GOL	D	305	-	5,5,5	0.89	0	5,5,5	1.04	0
5	PO4	E	301	-	4,4,4	0.93	0	6,6,6	0.41	0
6	GOL	B	302	-	5,5,5	0.91	0	5,5,5	1.00	0
5	PO4	D	302	-	4,4,4	0.91	0	6,6,6	0.45	0
6	GOL	A	303	-	5,5,5	0.92	0	5,5,5	0.96	0

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
6	GOL	D	304	-	-	3/4/4/4	-
6	GOL	A	302	-	-	2/4/4/4	-
6	GOL	D	303	-	-	2/4/4/4	-
6	GOL	D	305	-	-	2/4/4/4	-
6	GOL	B	302	-	-	4/4/4/4	-
6	GOL	A	303	-	-	2/4/4/4	-
6	GOL	B	301	-	-	0/4/4/4	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (15) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
6	A	302	GOL	O1-C1-C2-C3

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Mol	Chain	Res	Type	Atoms
6	B	302	GOL	C1-C2-C3-O3
6	B	302	GOL	O2-C2-C3-O3
6	D	303	GOL	O1-C1-C2-O2
6	D	304	GOL	C1-C2-C3-O3
6	A	303	GOL	O1-C1-C2-C3
6	D	303	GOL	O1-C1-C2-C3
6	A	303	GOL	O1-C1-C2-O2
6	A	302	GOL	O1-C1-C2-O2
6	D	305	GOL	O1-C1-C2-O2
6	B	302	GOL	O1-C1-C2-C3
6	D	304	GOL	O2-C2-C3-O3
6	B	302	GOL	O1-C1-C2-O2
6	D	304	GOL	O1-C1-C2-C3
6	D	305	GOL	O1-C1-C2-C3

There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	D	304	GOL	1	0
6	B	301	GOL	1	0

5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

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	A	220/220 (100%)	-0.20	2 (0%)	84	83	26, 43, 64, 126	0
1	D	219/220 (99%)	-0.17	1 (0%)	91	89	28, 45, 79, 98	0
2	B	223/225 (99%)	-0.14	1 (0%)	92	91	26, 42, 62, 132	0
2	E	223/225 (99%)	-0.15	2 (0%)	84	83	23, 43, 69, 118	0
3	C	8/18 (44%)	0.53	2 (25%)	0	0	57, 70, 90, 103	0
3	F	8/18 (44%)	1.10	2 (25%)	0	0	53, 76, 102, 111	0
All	All	901/926 (97%)	-0.15	10 (1%)	80	78	23, 43, 71, 132	0

All (10) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	141	THR	8.3
1	A	220	CYS	6.9
2	E	141	THR	4.3
3	F	75	VAL	3.1
3	C	75	VAL	2.7
3	F	78	ILE	2.6
2	E	142	ASN	2.6
1	A	87	GLU	2.1
3	C	82	ASP	2.1
1	D	208	THR	2.0

6.2 Non-standard residues in protein, DNA, RNA chains

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

6.3 Carbohydrates (i)

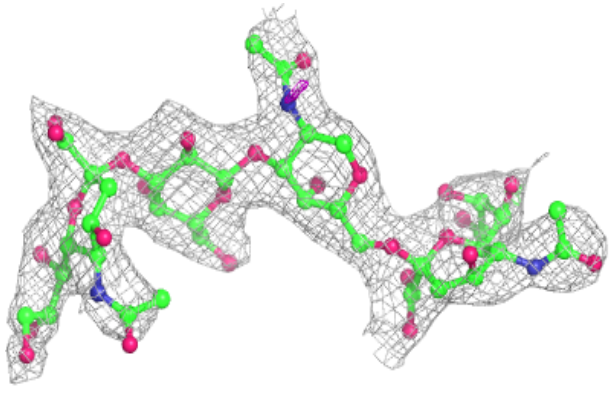
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95^{th} percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9'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

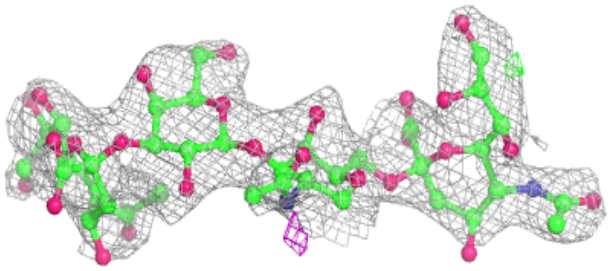
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
4	SIA	H	3	20/21	0.77	0.29	68,72,77,78	0
4	SIA	G	3	20/21	0.82	0.24	74,78,82,84	0
4	SIA	H	4	20/21	0.88	0.24	60,62,64,65	0
4	GAL	H	2	11/12	0.89	0.20	65,69,72,72	0
4	A2G	G	1	14/15	0.90	0.18	52,56,58,60	0
4	SIA	G	4	20/21	0.92	0.20	50,51,54,55	0
4	GAL	G	2	11/12	0.92	0.19	59,62,65,69	0
4	A2G	H	1	14/15	0.93	0.15	47,54,59,62	0

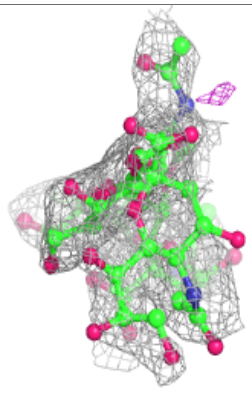
The following is a graphical depiction of the model fit to experimental electron density for oligosaccharide. Each fit is shown from different orientation to approximate a three-dimensional view.

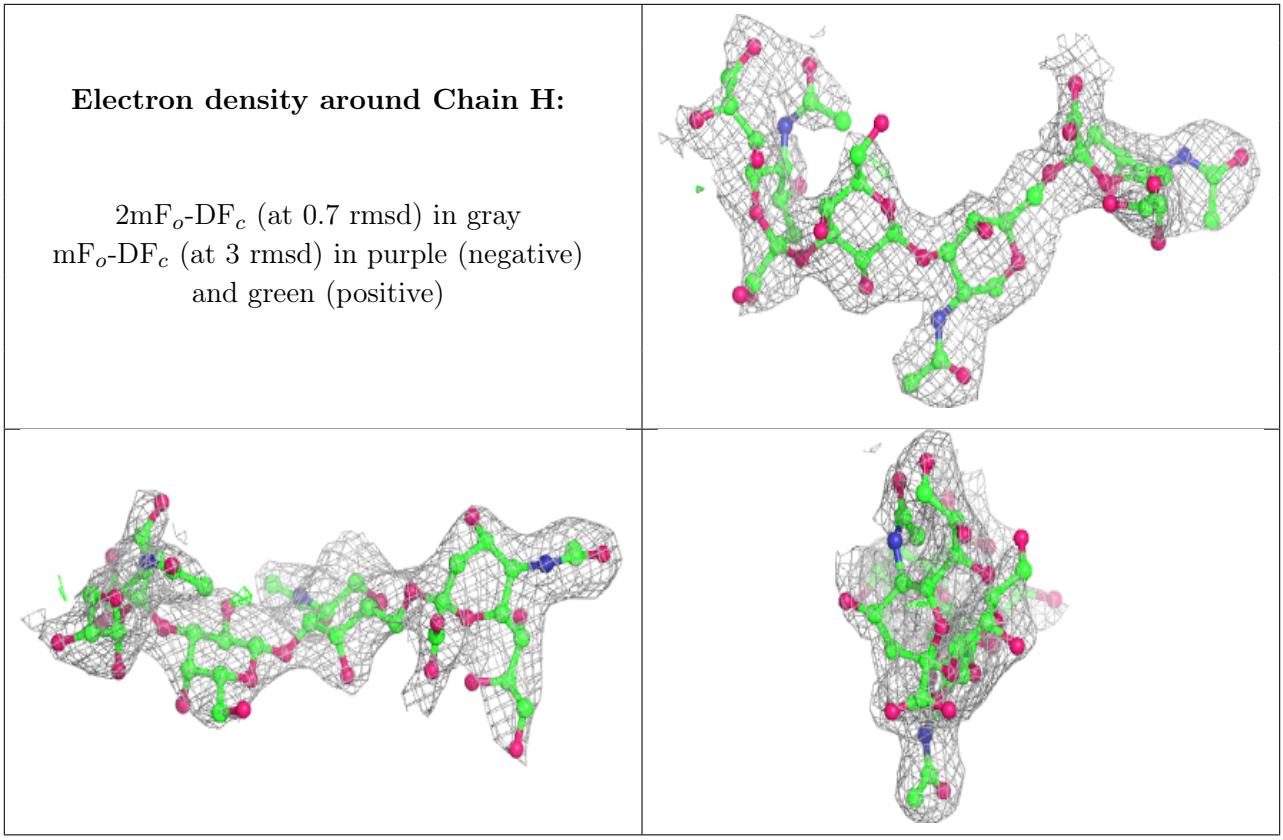
Electron density around Chain G:

$2mF_o$-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, 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
5	PO4	D	301	5/5	0.80	0.35	105,107,107,108	0
6	GOL	D	303	6/6	0.89	0.28	51,54,55,56	0
6	GOL	A	303	6/6	0.90	0.15	56,59,60,60	0
6	GOL	B	302	6/6	0.91	0.22	58,58,58,59	0
6	GOL	B	301	6/6	0.91	0.18	45,47,48,50	0
5	PO4	E	301	5/5	0.92	0.23	82,82,83,83	0
6	GOL	A	302	6/6	0.92	0.28	39,43,45,48	0
6	GOL	D	304	6/6	0.92	0.19	59,62,62,63	0
5	PO4	D	302	5/5	0.93	0.23	90,90,91,92	0
5	PO4	A	301	5/5	0.94	0.23	103,103,103,103	0
6	GOL	D	305	6/6	0.98	0.26	39,40,42,44	0

6.5 Other polymers (i)

There are no such residues in this entry.