



Full wwPDB X-ray Structure Validation Report ⓘ

Apr 11, 2024 – 03:29 PM EDT

PDB ID : 8TFH
Title : Ricin in complex with Fab JB4
Authors : Rudolph, M.J.; Mantis, N.
Deposited on : 2023-07-11
Resolution : 3.29 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 : 2.36.1
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) : 2.36.1

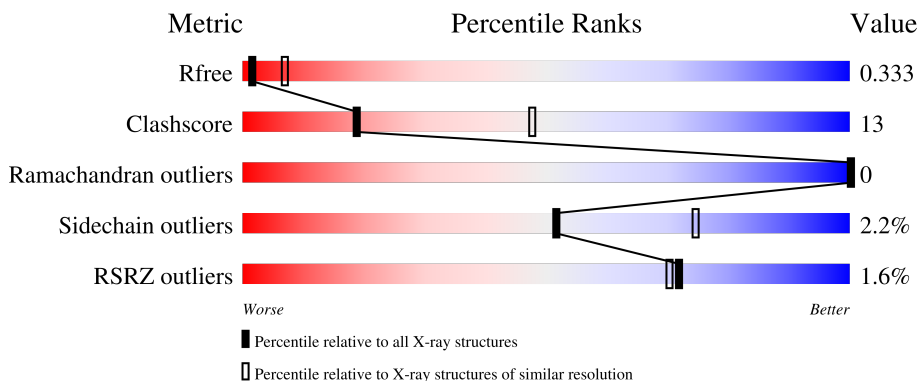
1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.29 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 $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	260	 3% 74% 25%
2	B	262	 0% 73% 26%
3	H	223	 0% 70% 28%
4	L	210	 2% 75% 24%
5	C	3	 67% 33%

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Mol	Chain	Length	Quality of chain
6	D	3	
6	E	3	

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
5	NAG	C	3	-	-	-	X
6	BMA	D	3	-	-	-	X

2 Entry composition [i](#)

There are 10 unique types of molecules in this entry. The entry contains 7598 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 Ricin A chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	260	2066	1308	368	385	5	0	1	0

- Molecule 2 is a protein called Ricin B chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	B	262	2035	1273	357	393	12	0	0	0

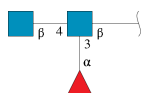
- Molecule 3 is a protein called JB4 monoclonal antibody heavy chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
3	H	223	1701	1077	283	334	7	0	0	0

- Molecule 4 is a protein called JB4 monoclonal antibody light chain.

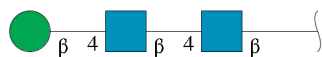
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
4	L	210	1637	1027	277	327	6	0	0	0

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



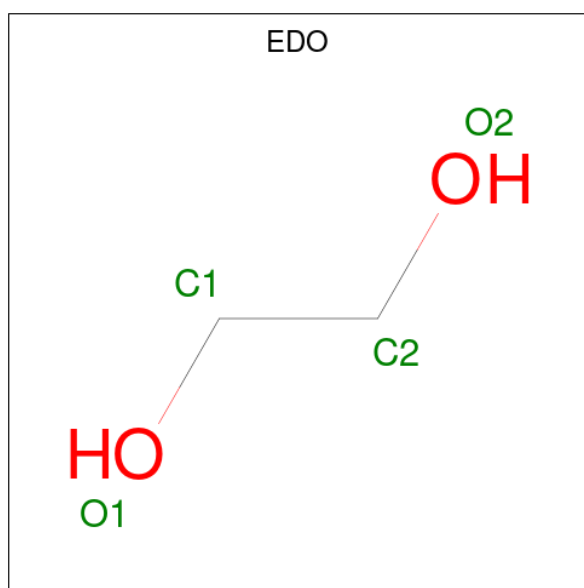
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
			Total	C	N	O			
5	C	3	38	22	2	14	0	0	0

- Molecule 6 is an oligosaccharide called 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	
6	D	3	Total	C	N	O	0	0	0
			39	22	2	15			
6	E	3	Total	C	N	O	0	0	0
			39	22	2	15			

- Molecule 7 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula: C₂H₆O₂).



Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	1	Total	C O	0	0
			4	2 2		
7	B	1	Total	C O	0	0
			4	2 2		
7	H	1	Total	C O	0	0
			4	2 2		

- Molecule 8 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf	
8	A	1	Total	Cl	0	0
			1	1		

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Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
8	L	2	Total Cl 2 2	0	0

- Molecule 9 is SULFATE ION (three-letter code: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
9	B	1	Total O S 5 4 1	0	0
9	H	1	Total O S 5 4 1	0	0

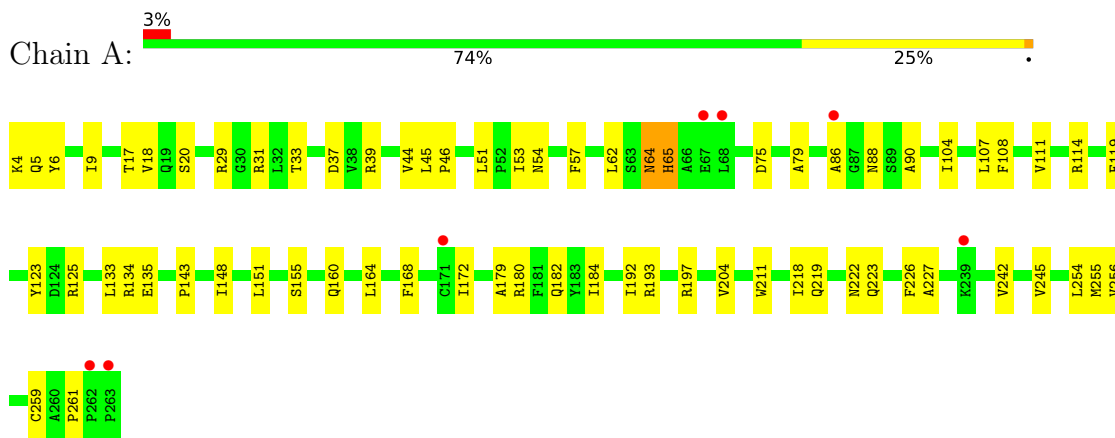
- Molecule 10 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
10	A	3	Total O 3 3	0	0
10	B	6	Total O 6 6	0	0
10	H	4	Total O 4 4	0	0
10	L	5	Total O 5 5	0	0

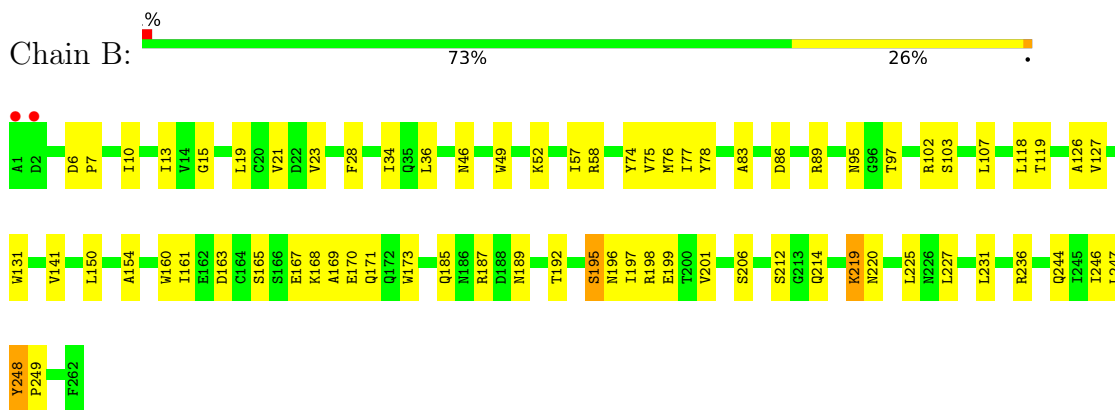
3 Residue-property plots

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.

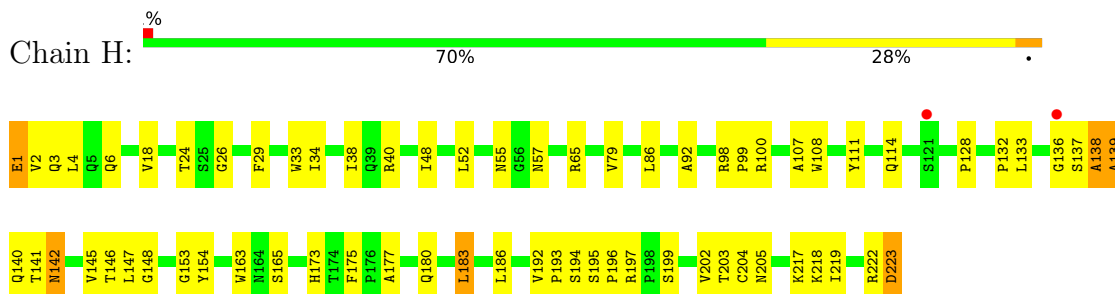
- Molecule 1: Ricin A chain



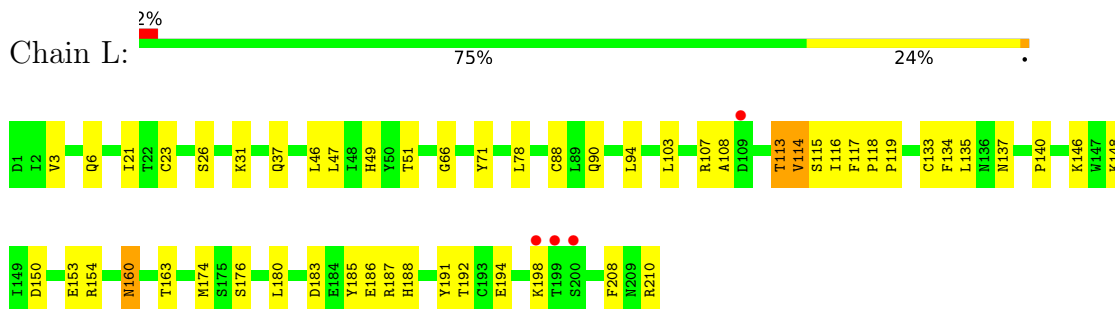
- Molecule 2: Ricin B chain



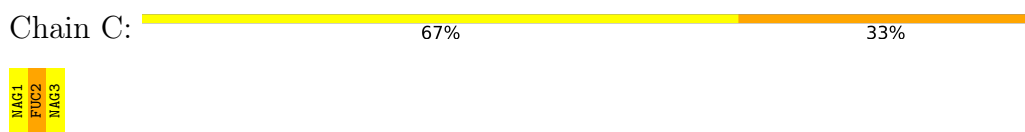
- Molecule 3: JB4 monoclonal antibody heavy chain



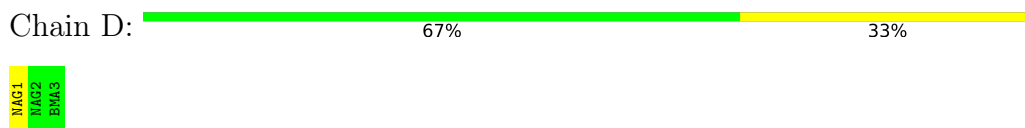
- Molecule 4: JB4 monoclonal antibody light chain



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



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



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



4 Data and refinement statistics

Property	Value	Source
Space group	I 41 2 2	Depositor
Cell constants a, b, c, α , β , γ	139.36Å 139.36Å 444.61Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	33.84 – 3.29 33.84 – 3.29	Depositor EDS
% Data completeness (in resolution range)	96.5 (33.84-3.29) 96.6 (33.84-3.29)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.57 (at 3.32Å)	Xtrriage
Refinement program	PHENIX (1.20.1_4487: ???)	Depositor
R, R_{free}	0.274 , 0.334 0.274 , 0.333	Depositor DCC
R_{free} test set	1614 reflections (4.94%)	wwPDB-VP
Wilson B-factor (Å ²)	64.1	Xtrriage
Anisotropy	0.074	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 52.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.28$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.85	EDS
Total number of atoms	7598	wwPDB-VP
Average B, all atoms (Å ²)	75.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.32% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: SO4, CL, NAG, FUC, EDO, BMA

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	0/2111	0.56	1/2872 (0.0%)
2	B	0.25	0/2080	0.54	1/2842 (0.0%)
3	H	0.28	0/1749	0.65	3/2393 (0.1%)
4	L	0.26	0/1676	0.53	1/2273 (0.0%)
All	All	0.26	0/7616	0.57	6/10380 (0.1%)

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
3	H	0	3
4	L	0	1
All	All	0	4

There are no bond length outliers.

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
2	B	195	SER	C-N-CA	7.97	141.62	121.70
1	A	64	ASN	C-N-CA	7.80	141.21	121.70
3	H	183	LEU	CA-CB-CG	6.86	131.07	115.30
4	L	114	VAL	C-N-CA	6.09	136.92	121.70
3	H	139	ALA	C-N-CA	6.08	136.89	121.70
3	H	183	LEU	CB-CG-CD2	5.39	120.17	111.00

There are no chirality outliers.

All (4) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
3	H	1	GLU	Peptide
3	H	138	ALA	Peptide
3	H	142	ASN	Peptide
4	L	113	THR	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	2066	0	2037	45	0
2	B	2035	0	1974	50	0
3	H	1701	0	1641	64	0
4	L	1637	0	1585	44	0
5	C	38	0	34	3	0
6	D	39	0	34	0	0
6	E	39	0	34	1	0
7	A	4	0	6	1	0
7	B	4	0	6	0	0
7	H	4	0	6	1	0
8	A	1	0	0	1	0
8	L	2	0	0	0	0
9	B	5	0	0	0	0
9	H	5	0	0	0	0
10	A	3	0	0	2	0
10	B	6	0	0	1	0
10	H	4	0	0	1	0
10	L	5	0	0	0	0
All	All	7598	0	7357	195	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 13.

All (195) 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)
3:H:147:LEU:HD13	3:H:219:ILE:HD12	1.36	1.04
3:H:99:PRO:HB3	3:H:107:ALA:HA	1.39	1.03

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:H:147:LEU:HD13	3:H:219:ILE:CD1	1.91	1.01
5:C:2:FUC:H5	5:C:3:NAG:H2	1.48	0.96
1:A:143:PRO:HD3	1:A:197[B]:ARG:HG2	1.52	0.90
1:A:193:ARG:NH2	8:A:302:CL:CL	2.44	0.88
3:H:147:LEU:CD1	3:H:219:ILE:HD12	2.06	0.85
4:L:114:VAL:HG21	4:L:133:CYS:HA	1.58	0.84
3:H:2:VAL:HG12	3:H:3:GLN:HB2	1.65	0.78
2:B:165:SER:HB3	2:B:168:LYS:HB2	1.64	0.78
2:B:231:LEU:HD23	2:B:247:LEU:HB3	1.67	0.74
2:B:236:ARG:NH1	2:B:244:GLN:OE1	2.19	0.74
1:A:222:ASN:HB2	2:B:6:ASP:HB2	1.71	0.72
3:H:4:LEU:HD22	3:H:24:THR:HG22	1.71	0.72
3:H:132:PRO:HD3	3:H:217:LYS:HD3	1.72	0.72
4:L:160:ASN:HB3	4:L:176:SER:HA	1.72	0.71
1:A:6:TYR:OH	1:A:53:ILE:O	2.08	0.70
2:B:219:LYS:HD2	2:B:225:LEU:HD12	1.73	0.70
1:A:46:PRO:HB2	1:A:51:LEU:HD11	1.73	0.70
3:H:222:ARG:NH1	4:L:119:PRO:O	2.25	0.70
4:L:107:ARG:HG2	4:L:108:ALA:H	1.58	0.69
1:A:44:VAL:HG12	1:A:255:MET:HB2	1.75	0.69
2:B:225:LEU:HD22	2:B:227:LEU:H	1.57	0.68
4:L:90:GLN:NE2	4:L:94:LEU:O	2.28	0.67
3:H:1:GLU:HG2	3:H:111:TYR:HE1	1.60	0.67
1:A:4:LYS:NZ	10:A:402:HOH:O	2.28	0.66
1:A:64:ASN:HB2	1:A:65:HIS:C	2.15	0.66
3:H:147:LEU:HD12	3:H:148:GLY:N	2.12	0.65
1:A:104:ILE:HA	1:A:107:LEU:HD23	1.78	0.65
3:H:6:GLN:H	3:H:114:GLN:HE22	1.43	0.65
3:H:203:THR:HG22	3:H:218:LYS:HG3	1.79	0.65
5:C:2:FUC:H5	5:C:3:NAG:C2	2.24	0.65
3:H:147:LEU:HD13	3:H:219:ILE:HD13	1.77	0.64
3:H:26:GLY:H	7:H:302:EDO:H11	1.63	0.64
3:H:38:ILE:HB	3:H:48:ILE:HD11	1.78	0.64
4:L:3:VAL:HG22	4:L:26:SER:HB3	1.80	0.63
2:B:197:ILE:HG22	2:B:198:ARG:H	1.64	0.62
3:H:196:PRO:HA	3:H:199:SER:OG	2.00	0.62
3:H:137:SER:OG	3:H:223:ASP:O	2.18	0.61
4:L:186:GLU:O	4:L:210:ARG:NH2	2.33	0.61
3:H:34:ILE:HG13	3:H:79:VAL:HG21	1.83	0.60
2:B:225:LEU:CD2	2:B:227:LEU:H	2.15	0.59
4:L:154:ARG:HH21	4:L:180:LEU:HD22	1.66	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:222:ASN:O	1:A:223:GLN:HG2	2.03	0.58
3:H:173:HIS:NE2	4:L:137:ASN:OD1	2.32	0.58
2:B:75:VAL:HG21	2:B:107:LEU:HD12	1.86	0.58
3:H:128:PRO:HB3	3:H:154:TYR:HB3	1.84	0.58
2:B:86:ASP:HA	2:B:89:ARG:HG2	1.85	0.57
3:H:193:PRO:O	3:H:196:PRO:HD2	2.05	0.57
1:A:227:ALA:HB2	2:B:7:PRO:HG3	1.87	0.57
4:L:114:VAL:HB	4:L:115:SER:C	2.26	0.56
4:L:114:VAL:HG22	4:L:134:PHE:N	2.20	0.56
4:L:37:GLN:HB2	4:L:47:LEU:HD11	1.86	0.56
1:A:133:LEU:HD13	7:A:301:EDO:H11	1.87	0.56
2:B:52:LYS:HG3	2:B:58:ARG:HD2	1.86	0.56
3:H:197:ARG:HE	3:H:197:ARG:C	2.07	0.56
3:H:1:GLU:HG2	3:H:111:TYR:CE1	2.39	0.56
1:A:182:GLN:NE2	2:B:220:ASN:O	2.39	0.56
2:B:13:ILE:HB	2:B:21:VAL:HB	1.88	0.55
1:A:64:ASN:OD1	1:A:148:ILE:HB	2.06	0.55
3:H:146:THR:OG1	4:L:114:VAL:O	2.23	0.55
3:H:55:ASN:ND2	3:H:57:ASN:OD1	2.40	0.55
3:H:139:ALA:HB3	3:H:140:GLN:HA	1.89	0.55
1:A:90:ALA:HB2	1:A:111:VAL:HG11	1.89	0.55
4:L:183:ASP:O	4:L:187:ARG:HD3	2.07	0.55
2:B:97:THR:OG1	10:B:401:HOH:O	2.17	0.54
4:L:146:LYS:NZ	4:L:153:GLU:HB3	2.22	0.54
3:H:147:LEU:HD11	3:H:202:VAL:HG11	1.90	0.54
3:H:136:GLY:HA2	3:H:222:ARG:HB2	1.90	0.54
1:A:222:ASN:ND2	2:B:6:ASP:OD2	2.41	0.53
2:B:195:SER:HB2	2:B:197:ILE:N	2.23	0.53
2:B:201:VAL:HG13	2:B:236:ARG:HH12	1.73	0.53
3:H:2:VAL:HB	3:H:26:GLY:HA3	1.91	0.53
1:A:4:LYS:O	1:A:5:GLN:NE2	2.31	0.53
3:H:100:ARG:HD2	3:H:108:TRP:CH2	2.44	0.52
3:H:192:VAL:HG22	3:H:193:PRO:HD2	1.92	0.52
3:H:197:ARG:O	3:H:197:ARG:NE	2.32	0.52
1:A:45:LEU:HD21	1:A:254:LEU:HD13	1.92	0.52
4:L:114:VAL:HG22	4:L:134:PHE:O	2.10	0.52
2:B:126:ALA:HA	2:B:212:SER:HA	1.92	0.51
3:H:141:THR:HG23	3:H:142:ASN:HB2	1.93	0.51
2:B:192:THR:HG23	2:B:214:GLN:HG2	1.92	0.51
3:H:165:SER:N	3:H:205:ASN:OD1	2.27	0.51
3:H:137:SER:HA	3:H:138:ALA:C	2.31	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:L:114:VAL:HB	4:L:116:ILE:N	2.26	0.51
3:H:177:ALA:HB2	3:H:186:LEU:HD23	1.93	0.50
3:H:177:ALA:HA	3:H:186:LEU:HB3	1.92	0.50
4:L:114:VAL:CG1	4:L:117:PHE:CE2	2.95	0.50
4:L:146:LYS:HZ1	4:L:153:GLU:HB3	1.75	0.50
1:A:29:ARG:NH1	1:A:179:ALA:O	2.40	0.50
4:L:113:THR:HG22	4:L:114:VAL:H	1.76	0.50
3:H:140:GLN:HG2	3:H:141:THR:H	1.74	0.50
1:A:86:ALA:HB3	1:A:155:SER:HB3	1.94	0.49
2:B:74:TYR:CE2	2:B:76:MET:HB3	2.46	0.49
2:B:127:VAL:HG21	2:B:214:GLN:HB2	1.94	0.49
2:B:154:ALA:HB2	2:B:173:TRP:CH2	2.47	0.49
5:C:2:FUC:C5	5:C:3:NAG:H2	2.33	0.49
4:L:31:LYS:O	4:L:51:THR:OG1	2.31	0.49
1:A:218:ILE:HG12	1:A:245:VAL:HG12	1.93	0.49
2:B:34:ILE:HD11	2:B:77:ILE:HB	1.93	0.49
2:B:103:SER:HB2	3:H:33:TRP:HZ2	1.78	0.49
3:H:193:PRO:HB2	3:H:196:PRO:HD3	1.94	0.49
4:L:135:LEU:N	4:L:174:MET:O	2.31	0.49
1:A:119:PHE:HB2	1:A:125:ARG:HG2	1.95	0.49
2:B:74:TYR:HA	2:B:119:THR:HG22	1.94	0.48
3:H:18:VAL:HG12	3:H:86:LEU:HD11	1.93	0.48
1:A:123:TYR:CD1	1:A:134:ARG:HD3	2.48	0.48
1:A:261:PRO:O	10:A:401:HOH:O	2.20	0.48
2:B:195:SER:HB2	2:B:197:ILE:HG13	1.96	0.48
2:B:150:LEU:HB2	2:B:161:ILE:HG23	1.95	0.48
2:B:163:ASP:N	2:B:163:ASP:OD1	2.47	0.47
3:H:145:VAL:HG23	3:H:194:SER:HB3	1.97	0.47
1:A:108:PHE:HB3	1:A:111:VAL:HG13	1.97	0.47
2:B:23:VAL:HB	2:B:46:ASN:HB2	1.97	0.47
4:L:114:VAL:HG12	4:L:117:PHE:CE2	2.50	0.47
4:L:118:PRO:HB3	4:L:208:PHE:CE1	2.50	0.47
1:A:17:THR:HG23	1:A:20:SER:H	1.80	0.47
2:B:78:TYR:CD2	2:B:83:ALA:HB2	2.50	0.47
3:H:193:PRO:C	3:H:196:PRO:HD2	2.34	0.47
2:B:95:ASN:HB2	2:B:227:LEU:HD11	1.96	0.47
3:H:141:THR:HA	3:H:142:ASN:HA	1.54	0.47
3:H:98:ARG:NH2	3:H:100:ARG:O	2.47	0.46
4:L:148:LYS:HB2	4:L:192:THR:HG22	1.96	0.46
1:A:168:PHE:O	1:A:172:ILE:HG13	2.15	0.46
4:L:133:CYS:N	4:L:176:SER:O	2.41	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:H:180:GLN:N	3:H:183:LEU:O	2.48	0.46
1:A:33:THR:HG22	1:A:45:LEU:HA	1.98	0.46
1:A:18:VAL:HG12	1:A:192:ILE:HG22	1.97	0.46
3:H:24:THR:OG1	3:H:29:PHE:CD1	2.69	0.46
2:B:197:ILE:HG22	2:B:198:ARG:N	2.30	0.46
3:H:133:LEU:HD13	4:L:117:PHE:CE2	2.50	0.46
4:L:150:ASP:HB2	4:L:188:HIS:HD2	1.80	0.46
4:L:185:TYR:O	4:L:191:TYR:OH	2.31	0.45
4:L:66:GLY:HA3	4:L:71:TYR:CD1	2.51	0.45
1:A:75:ASP:O	1:A:79:ALA:N	2.42	0.45
2:B:199:GLU:O	2:B:246:ILE:HB	2.17	0.45
1:A:226:PHE:HE2	1:A:242:VAL:HG12	1.81	0.45
3:H:203:THR:CG2	3:H:218:LYS:HG3	2.46	0.45
1:A:184:ILE:HG12	1:A:204:VAL:HG13	1.99	0.45
4:L:113:THR:HG22	4:L:114:VAL:HG23	1.98	0.44
4:L:46:LEU:HD21	4:L:49:HIS:HB3	2.00	0.44
3:H:148:GLY:HA2	3:H:163:TRP:CZ2	2.52	0.44
3:H:163:TRP:CD1	3:H:202:VAL:HG12	2.52	0.44
3:H:33:TRP:CE2	3:H:52:LEU:HD22	2.53	0.44
1:A:53:ILE:HD12	1:A:54:ASN:N	2.33	0.43
1:A:9:ILE:HD11	1:A:31:ARG:HG3	2.00	0.43
2:B:168:LYS:HD3	2:B:169:ALA:H	1.83	0.43
3:H:40:ARG:HB3	3:H:92:ALA:HB2	2.00	0.43
3:H:141:THR:CG2	3:H:142:ASN:HB2	2.48	0.43
2:B:15:GLY:HA3	2:B:131:TRP:CE3	2.54	0.43
4:L:146:LYS:HB3	4:L:194:GLU:HB3	2.00	0.43
3:H:141:THR:HG23	3:H:142:ASN:CB	2.48	0.43
3:H:24:THR:OG1	3:H:29:PHE:HD1	2.01	0.43
1:A:37:ASP:HB3	1:A:44:VAL:HG23	2.01	0.42
1:A:180:ARG:HA	1:A:254:LEU:HB2	2.01	0.42
2:B:169:ALA:HA	2:B:170:GLU:HA	1.76	0.42
3:H:147:LEU:HD12	3:H:147:LEU:C	2.39	0.42
4:L:6:GLN:HE21	4:L:21:ILE:HG21	1.84	0.42
2:B:23:VAL:HG21	2:B:28:PHE:HE1	1.83	0.42
2:B:189:ASN:HD22	2:B:206:SER:HA	1.83	0.42
2:B:225:LEU:HD22	2:B:227:LEU:N	2.31	0.42
4:L:187:ARG:O	4:L:188:HIS:CG	2.73	0.42
1:A:211:TRP:CH2	1:A:256:VAL:HB	2.54	0.42
2:B:36:LEU:HD13	2:B:118:LEU:HD11	2.00	0.42
4:L:78:LEU:HD11	4:L:103:LEU:HD21	2.00	0.42
1:A:39:ARG:NH1	1:A:219:GLN:OE1	2.53	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:143:PRO:HD3	1:A:197[A]:ARG:HG2	2.02	0.42
1:A:62:LEU:HD23	1:A:148:ILE:HD11	2.02	0.42
1:A:53:ILE:H	1:A:53:ILE:HG13	1.68	0.41
3:H:153:GLY:HA2	3:H:183:LEU:HD12	2.01	0.41
4:L:187:ARG:O	4:L:188:HIS:ND1	2.52	0.41
2:B:19:LEU:HB2	2:B:36:LEU:HG	2.01	0.41
2:B:195:SER:HA	2:B:196:ASN:HB2	2.02	0.41
3:H:33:TRP:CE3	3:H:52:LEU:HB2	2.55	0.41
3:H:147:LEU:CD1	3:H:219:ILE:CD1	2.76	0.41
4:L:94:LEU:HA	4:L:94:LEU:HD12	1.59	0.41
2:B:49:TRP:CE3	2:B:57:ILE:HG22	2.55	0.41
2:B:170:GLU:HG2	2:B:185:GLN:HG2	2.02	0.41
3:H:193:PRO:HB2	3:H:196:PRO:CD	2.50	0.41
1:A:135:GLU:H	1:A:135:GLU:CD	2.24	0.41
3:H:133:LEU:HD12	3:H:133:LEU:O	2.20	0.41
2:B:10:ILE:HG21	6:E:2:NAG:H83	2.03	0.41
4:L:117:PHE:HE2	4:L:134:PHE:CE2	2.39	0.41
2:B:195:SER:HB2	2:B:196:ASN:C	2.40	0.41
2:B:248:TYR:CG	2:B:249:PRO:HD2	2.56	0.41
1:A:151:LEU:HD12	1:A:164:LEU:HD22	2.03	0.41
3:H:175:PHE:CD1	4:L:163:THR:HG23	2.56	0.41
10:H:401:HOH:O	4:L:94:LEU:HD11	2.20	0.41
2:B:10:ILE:HD12	2:B:10:ILE:H	1.85	0.41
3:H:133:LEU:HD13	4:L:117:PHE:CD2	2.57	0.41
1:A:6:TYR:CE2	1:A:57:PHE:HB2	2.55	0.40
1:A:160:GLN:H	1:A:160:GLN:HG3	1.67	0.40
4:L:107:ARG:HG2	4:L:108:ALA:N	2.33	0.40
2:B:141:VAL:HA	2:B:173:TRP:O	2.21	0.40
4:L:140:PRO:HG2	4:L:198:LYS:HE2	2.04	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	259/260 (100%)	246 (95%)	13 (5%)	0	100	100
2	B	260/262 (99%)	242 (93%)	18 (7%)	0	100	100
3	H	221/223 (99%)	201 (91%)	20 (9%)	0	100	100
4	L	208/210 (99%)	198 (95%)	10 (5%)	0	100	100
All	All	948/955 (99%)	887 (94%)	61 (6%)	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	220/219 (100%)	216 (98%)	4 (2%)	59	78
2	B	227/227 (100%)	220 (97%)	7 (3%)	40	67
3	H	190/190 (100%)	186 (98%)	4 (2%)	53	75
4	L	186/186 (100%)	183 (98%)	3 (2%)	62	79
All	All	823/822 (100%)	805 (98%)	18 (2%)	52	74

All (18) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	65	HIS
1	A	88	ASN
1	A	114	ARG
1	A	259	CYS
2	B	102	ARG
2	B	160	TRP
2	B	167	GLU
2	B	171	GLN
2	B	187	ARG
2	B	219	LYS
2	B	248	TYR
3	H	65	ARG
3	H	195	SER

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Mol	Chain	Res	Type
3	H	204	CYS
3	H	223	ASP
4	L	23	CYS
4	L	88	CYS
4	L	160	ASN

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

Mol	Chain	Res	Type
3	H	142	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](#)

9 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
5	NAG	C	1	5,1	14,14,15	0.68	0	17,19,21	1.38	3 (17%)
5	FUC	C	2	5	10,10,11	1.28	1 (10%)	14,14,16	1.70	2 (14%)
5	NAG	C	3	5	14,14,15	0.32	0	17,19,21	0.74	0
6	NAG	D	1	6,2	14,14,15	0.61	0	17,19,21	1.73	4 (23%)
6	NAG	D	2	6	14,14,15	0.25	0	17,19,21	0.68	0
6	BMA	D	3	6	11,11,12	0.67	0	15,15,17	0.76	0
6	NAG	E	1	6,2	14,14,15	0.66	0	17,19,21	1.16	3 (17%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
6	NAG	E	2	6	14,14,15	0.26	0	17,19,21	0.46	0
6	BMA	E	3	6	11,11,12	0.58	0	15,15,17	0.77	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
5	NAG	C	1	5,1	-	1/6/23/26	0/1/1/1
5	FUC	C	2	5	-	-	0/1/1/1
5	NAG	C	3	5	-	2/6/23/26	0/1/1/1
6	NAG	D	1	6,2	-	1/6/23/26	0/1/1/1
6	NAG	D	2	6	-	2/6/23/26	0/1/1/1
6	BMA	D	3	6	-	2/2/19/22	0/1/1/1
6	NAG	E	1	6,2	-	2/6/23/26	0/1/1/1
6	NAG	E	2	6	-	4/6/23/26	0/1/1/1
6	BMA	E	3	6	-	2/2/19/22	0/1/1/1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	C	2	FUC	C4-C5	3.24	1.60	1.52

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	C	2	FUC	O5-C5-C4	4.38	117.38	109.52
6	D	1	NAG	C1-O5-C5	3.86	117.42	112.19
6	D	1	NAG	O4-C4-C3	-3.40	102.49	110.35
5	C	2	FUC	C1-O5-C5	3.20	120.03	112.78
5	C	1	NAG	C1-O5-C5	3.19	116.51	112.19
6	D	1	NAG	O5-C5-C6	-3.09	102.36	107.20
6	E	1	NAG	C1-C2-N2	2.73	115.16	110.49
6	D	1	NAG	C1-C2-N2	2.61	114.95	110.49
5	C	1	NAG	O5-C5-C6	-2.53	103.25	107.20
6	E	1	NAG	C1-O5-C5	2.23	115.21	112.19
6	E	1	NAG	O4-C4-C3	-2.20	105.25	110.35
5	C	1	NAG	O4-C4-C3	-2.13	105.43	110.35

There are no chirality outliers.

All (16) torsion outliers are listed below:

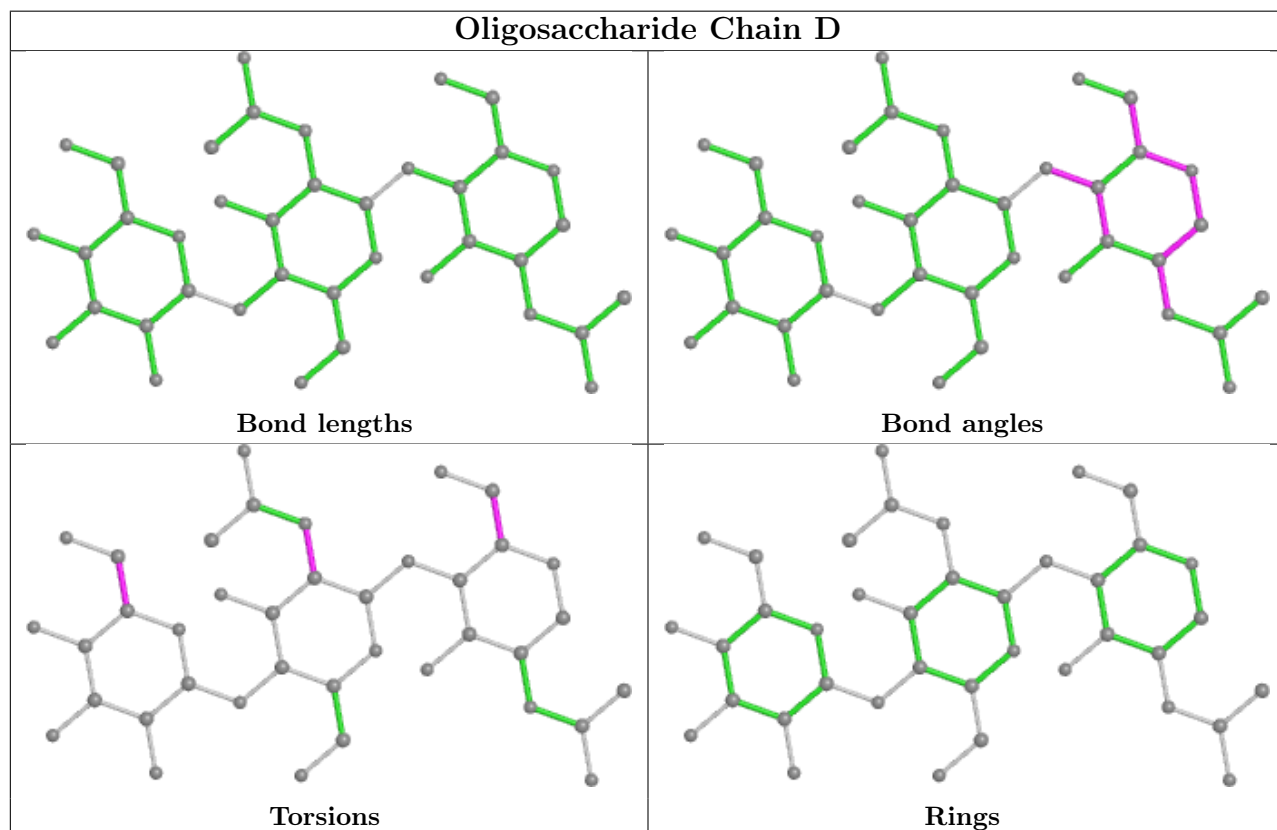
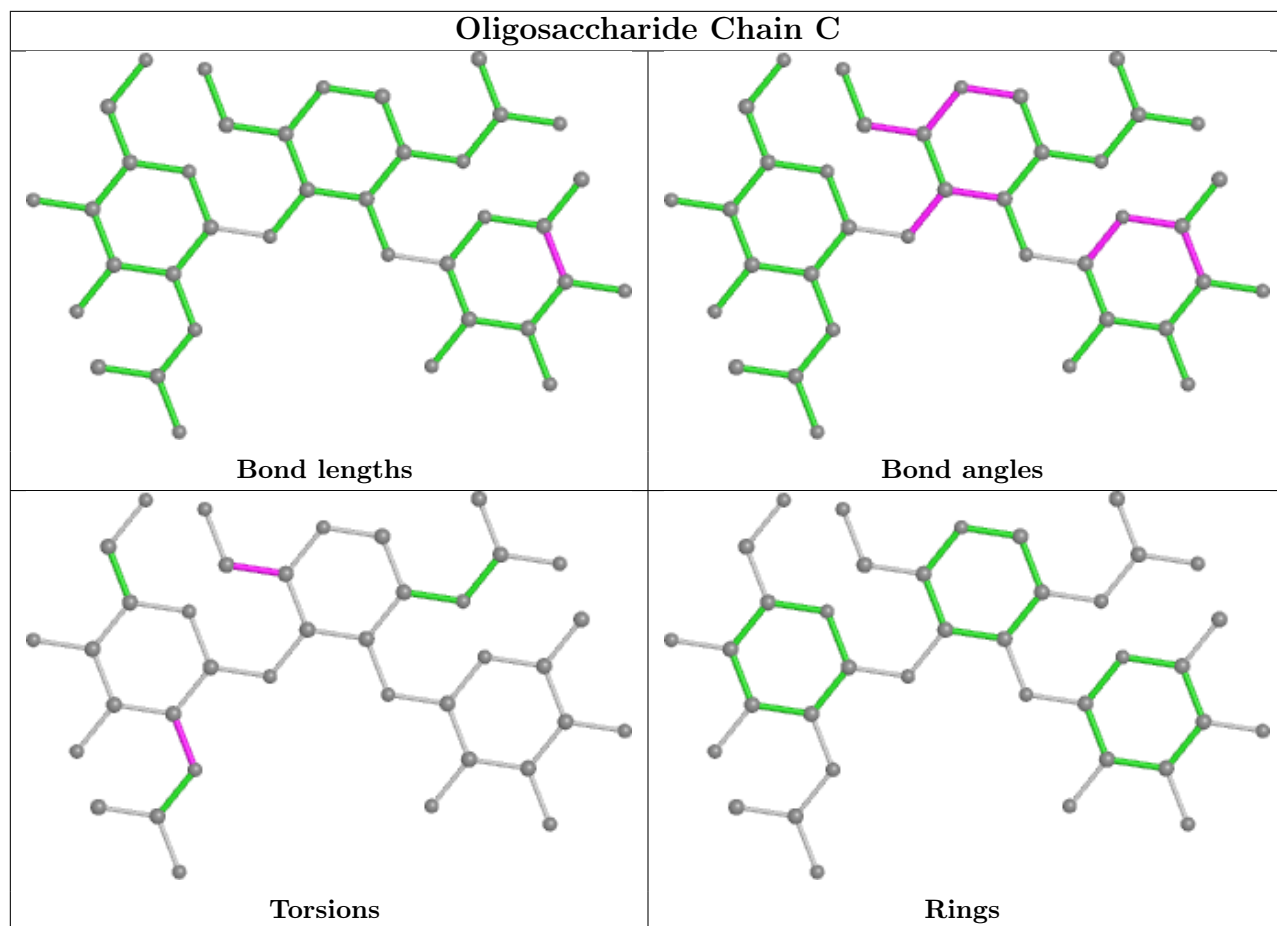
Mol	Chain	Res	Type	Atoms
6	E	1	NAG	O5-C5-C6-O6
6	E	3	BMA	O5-C5-C6-O6
6	E	2	NAG	C8-C7-N2-C2
6	E	2	NAG	O7-C7-N2-C2
6	E	1	NAG	C4-C5-C6-O6
6	E	2	NAG	O5-C5-C6-O6
6	E	2	NAG	C4-C5-C6-O6
5	C	1	NAG	O5-C5-C6-O6
6	D	1	NAG	O5-C5-C6-O6
6	E	3	BMA	C4-C5-C6-O6
6	D	3	BMA	C4-C5-C6-O6
6	D	3	BMA	O5-C5-C6-O6
5	C	3	NAG	C3-C2-N2-C7
6	D	2	NAG	C3-C2-N2-C7
5	C	3	NAG	C1-C2-N2-C7
6	D	2	NAG	C1-C2-N2-C7

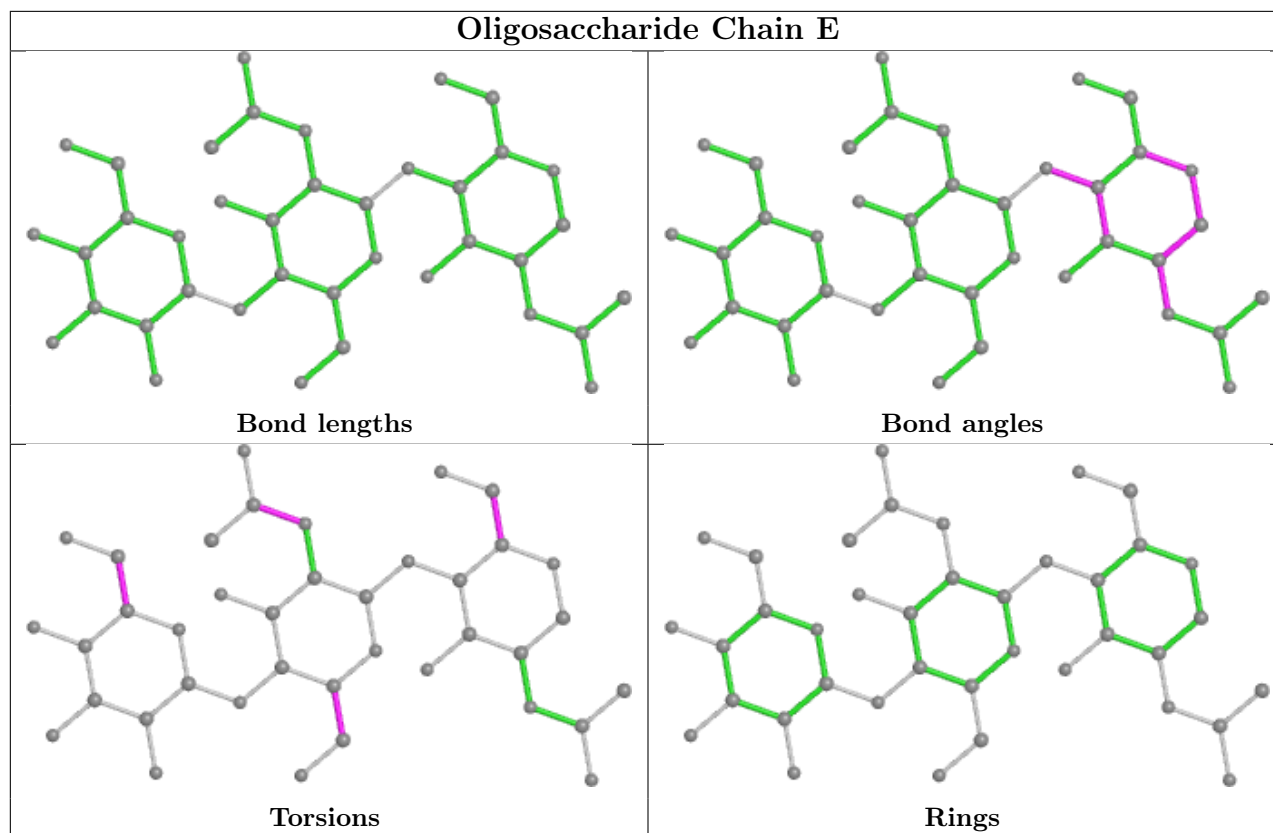
There are no ring outliers.

3 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	E	2	NAG	1	0
5	C	3	NAG	3	0
5	C	2	FUC	3	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](#)

Of 8 ligands modelled in this entry, 3 are monoatomic - leaving 5 for Mogul analysis.

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	SO4	B	301	-	4,4,4	0.14	0	6,6,6	0.07	0
7	EDO	A	301	-	3,3,3	0.43	0	2,2,2	0.36	0
7	EDO	H	302	-	3,3,3	0.43	0	2,2,2	0.40	0
9	SO4	H	301	-	4,4,4	0.13	0	6,6,6	0.11	0
7	EDO	B	302	-	3,3,3	0.46	0	2,2,2	0.34	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns.

'-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
7	EDO	A	301	-	-	1/1/1/1	-
7	EDO	H	302	-	-	0/1/1/1	-
7	EDO	B	302	-	-	0/1/1/1	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (1) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
7	A	301	EDO	O1-C1-C2-O2

There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
7	A	301	EDO	1	0
7	H	302	EDO	1	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data [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	260/260 (100%)	0.16	7 (2%) 54 52	40, 82, 122, 170	0
2	B	262/262 (100%)	-0.18	2 (0%) 86 86	38, 57, 86, 146	0
3	H	223/223 (100%)	0.11	2 (0%) 84 84	36, 75, 122, 160	0
4	L	210/210 (100%)	0.14	4 (1%) 66 65	44, 81, 114, 159	0
All	All	955/955 (100%)	0.05	15 (1%) 72 70	36, 73, 117, 170	0

All (15) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	263	PRO	3.9
2	B	1	ALA	3.3
1	A	68	LEU	3.1
1	A	239	LYS	2.9
1	A	262	PRO	2.9
4	L	199	THR	2.8
1	A	67	GLU	2.4
3	H	136	GLY	2.3
4	L	198	LYS	2.3
3	H	121	SER	2.3
1	A	171	CYS	2.3
1	A	86	ALA	2.2
4	L	200	SER	2.1
2	B	2	ASP	2.1
4	L	109	ASP	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

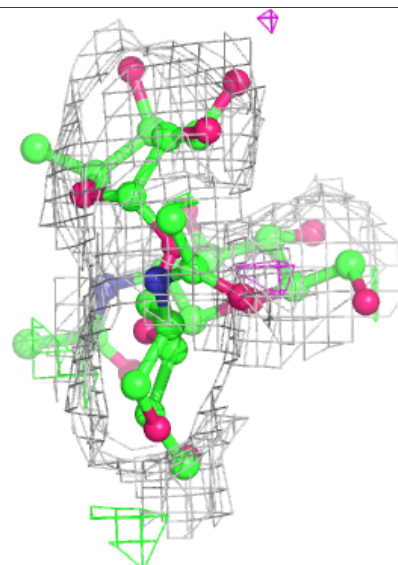
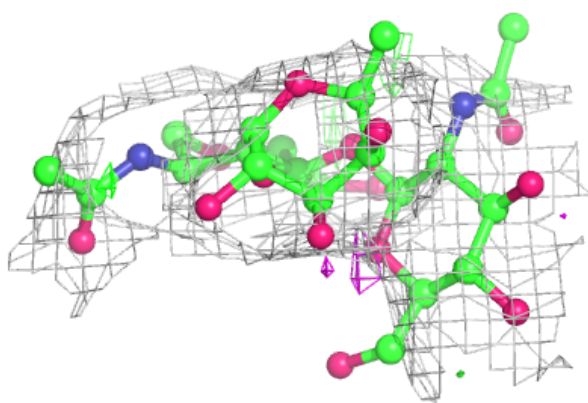
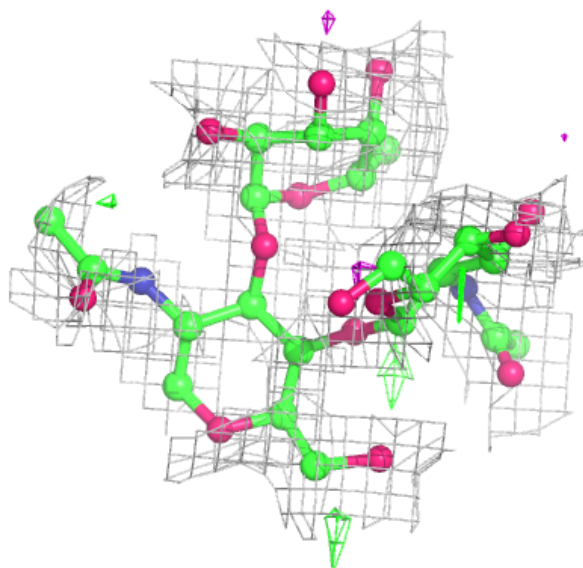
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q<0.9' lists 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	BMA	D	3	11/12	0.61	0.47	112,136,163,164	0
5	NAG	C	3	14/15	0.67	0.46	100,127,153,161	0
6	BMA	E	3	11/12	0.81	0.21	89,106,120,125	0
5	NAG	C	1	14/15	0.82	0.21	85,108,125,127	0
5	FUC	C	2	10/11	0.88	0.33	104,112,123,124	0
6	NAG	E	1	14/15	0.92	0.21	47,60,64,80	0
6	NAG	E	2	14/15	0.93	0.25	50,66,98,99	0
6	NAG	D	1	14/15	0.94	0.21	42,55,74,83	0
6	NAG	D	2	14/15	0.94	0.20	59,84,112,114	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.

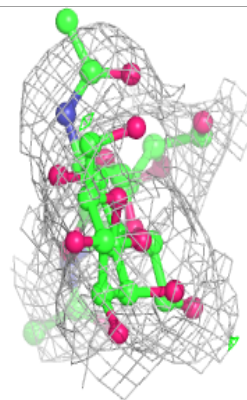
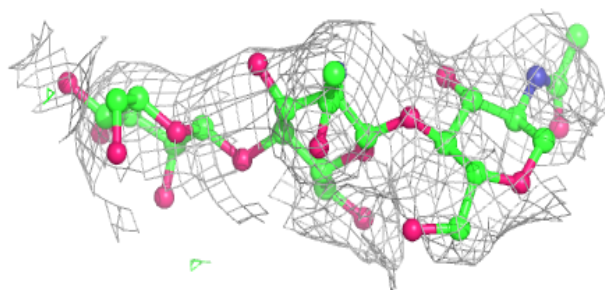
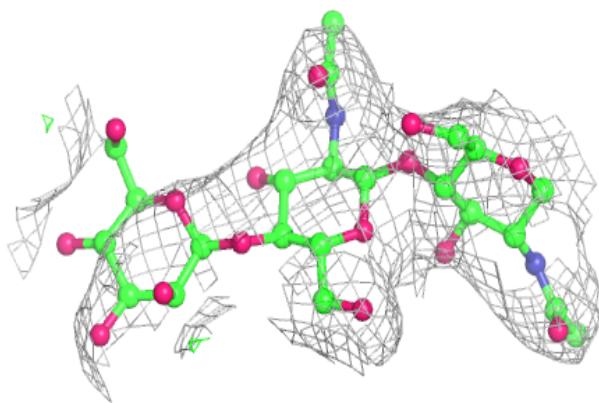
Electron density around Chain C:

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

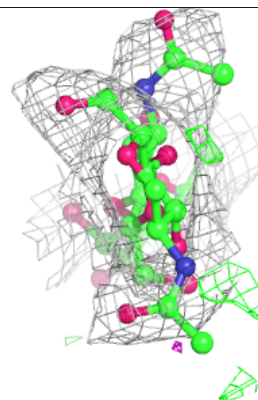
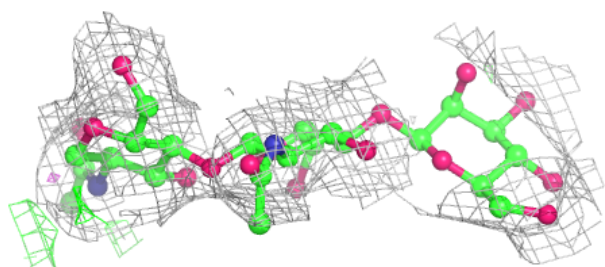
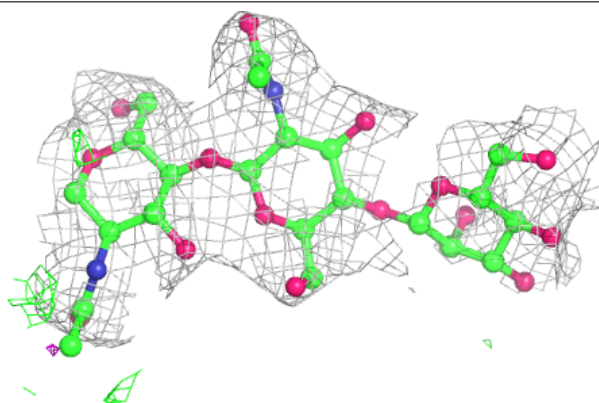


Electron density around Chain D:

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

**Electron density around Chain E:**

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



6.4 Ligands [i](#)

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
7	EDO	A	301	4/4	0.78	0.32	89,93,101,103	0
8	CL	A	302	1/1	0.86	0.14	61,61,61,61	0
8	CL	L	301	1/1	0.89	0.11	62,62,62,62	0
8	CL	L	302	1/1	0.91	0.10	55,55,55,55	0
7	EDO	H	302	4/4	0.92	0.18	24,46,53,57	0
9	SO4	B	301	5/5	0.94	0.22	80,87,89,107	0
9	SO4	H	301	5/5	0.94	0.21	84,93,98,108	0
7	EDO	B	302	4/4	0.95	0.11	28,35,38,43	0

6.5 Other polymers [i](#)

There are no such residues in this entry.