



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 20, 2026 – 09:07 AM UTC

PDB ID : 5E8E / pdb_00005e8e
Title : Crystal structure of thrombin bound to an exosite 1-specific IgA Fab
Authors : Baglin, T.P.; Langdown, J.; Frasson, R.; Huntington, J.A.
Deposited on : 2015-10-14
Resolution : 1.90 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

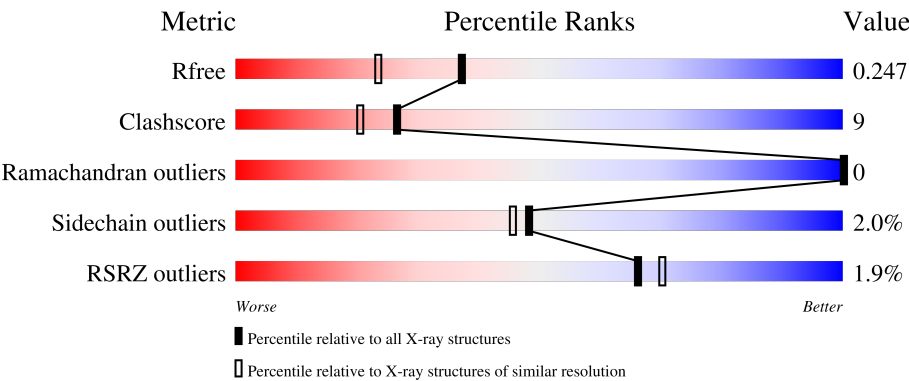
1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.90 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R _{free}	180053	7789 (1.90-1.90)
Clashscore	190562	8410 (1.90-1.90)
Ramachandran outliers	187476	8333 (1.90-1.90)
Sidechain outliers	187428	8333 (1.90-1.90)
RSRZ outliers	180081	7790 (1.90-1.90)

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	215	% 83%16%.
2	B	229	3% 78%21%.
3	H	259	% 89%11%
4	L	36	3% 72%14%6%8%
5	C	2	50%50%

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Mol	Chain	Length	Quality of chain
6	D	7	

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	2	-	-	X	-
8	PO4	B	301	-	-	X	-

2 Entry composition

There are 12 unique types of molecules in this entry. The entry contains 6373 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 IGA FAB LIGHT CHAIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	215	Total	C	N	O	S	0	5	0
			1670	1048	280	337	5			

- Molecule 2 is a protein called IGA FAB HEAVY CHAIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	229	Total	C	N	O	S	0	3	0
			1722	1088	283	340	11			

- Molecule 3 is a protein called THROMBIN HEAVY CHAIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	H	259	Total	C	N	O	S	0	1	0
			2091	1332	370	375	14			

- Molecule 4 is a protein called THROMBIN LIGHT CHAIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	L	33	Total	C	N	O	S	0	1	0
			270	169	42	58	1			

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



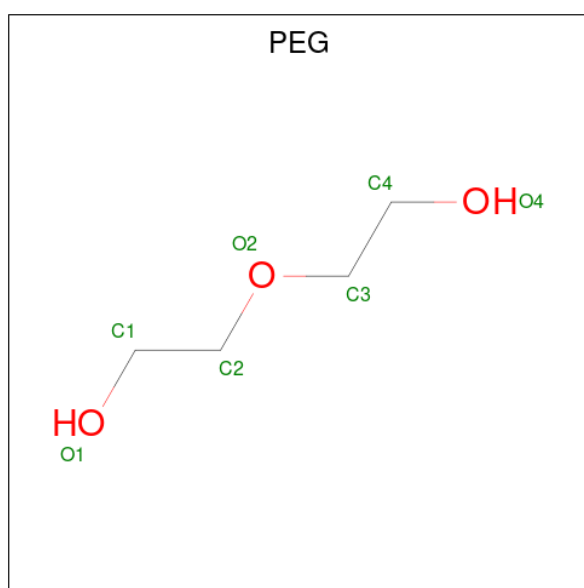
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
5	C	2	Total	C	N	O		0	0	0
			28	16	2	10				

- Molecule 6 is an oligosaccharide called N-acetyl-alpha-neuraminic acid-(2-6)-beta-D-galactopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-2)-alpha-D-mannopyranose-(1-6)-beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
6	D	7	Total	C	N	O	0	0	0
			95	53	4	38			

- Molecule 7 is DI(HYDROXYETHYL)ETHER (CCD ID: PEG) (formula: $C_4H_{10}O_3$).



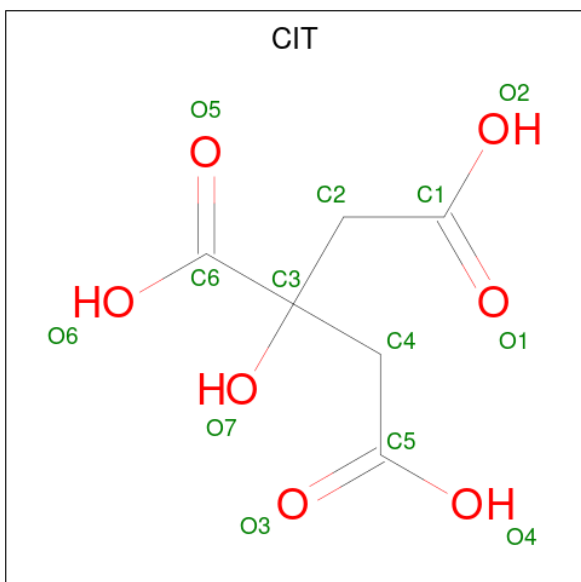
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
7	A	1	Total	C	O	0	0
			7	4	3		
7	H	1	Total	C	O	0	0
			7	4	3		
7	H	1	Total	C	O	0	0
			7	4	3		

- Molecule 8 is PHOSPHATE ION (CCD ID: PO4) (formula: O_4P).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
8	B	1	Total	O	P	0	0
			5	4	1		
8	H	1	Total	O	P	0	0
			5	4	1		
8	H	1	Total	O	P	0	0
			5	4	1		
8	H	1	Total	O	P	0	0
			5	4	1		
8	H	1	Total	O	P	0	0
			5	4	1		

- Molecule 9 is CITRIC ACID (CCD ID: CIT) (formula: $C_6H_8O_7$).

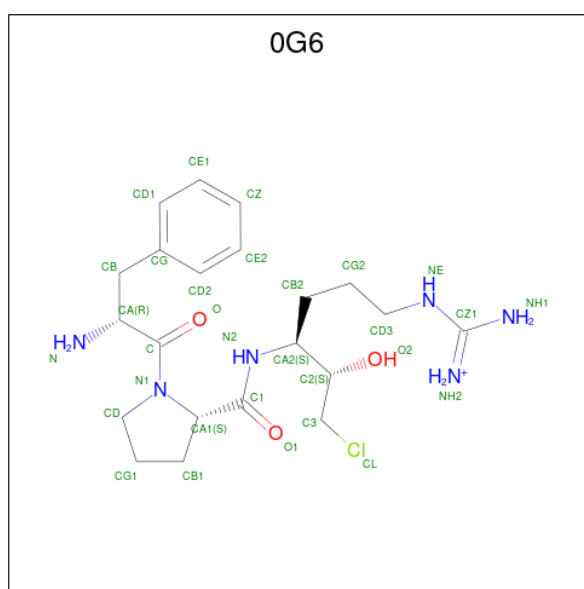


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
9	H	1	Total	C	O	0	0
			13	6	7		

- Molecule 10 is SODIUM ION (CCD ID: NA) (formula: Na).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
10	H	1	Total	Na	0	0
			1	1		

- Molecule 11 is D-phenylalanyl-N-[(2S,3S)-6-{{[amino(iminio)methyl]amino}-1-chloro-2-hydroxyhexan-3-yl]-L-prolinamide (CCD ID: 0G6) (formula: C₂₁H₃₄ClN₆O₃).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
11	H	1	Total	C	N	O	0	0
			30	21	6	3		

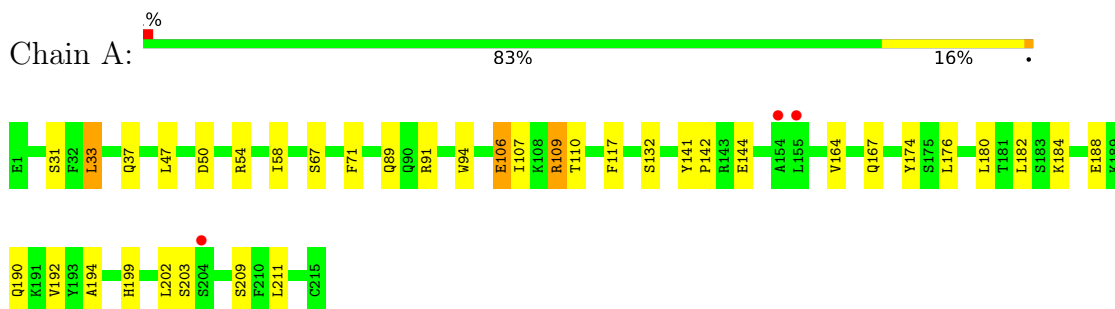
- Molecule 12 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
12	A	95	Total	O	0	0
			95	95		
12	B	105	Total	O	0	0
			105	105		
12	H	189	Total	O	0	0
			189	189		
12	L	18	Total	O	0	0
			18	18		

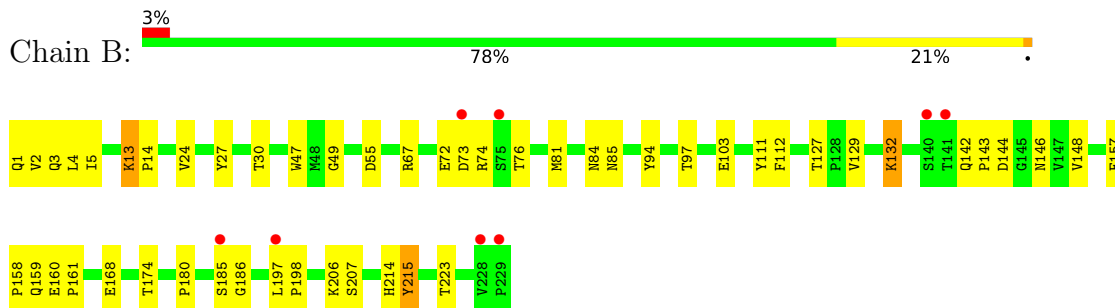
3 Residue-property plots [i](#)

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.

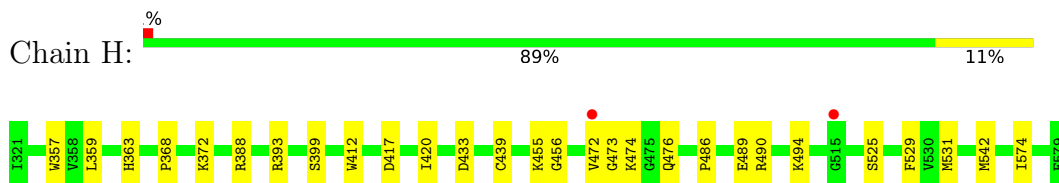
- Molecule 1: IGA FAB LIGHT CHAIN



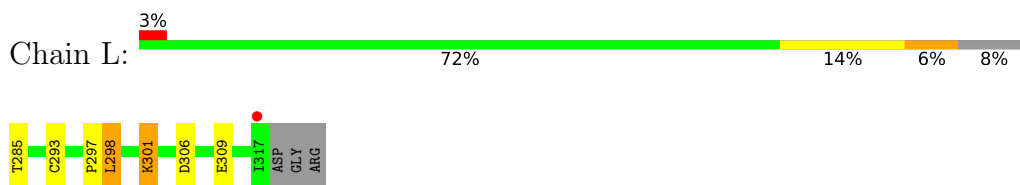
- Molecule 2: IGA FAB HEAVY CHAIN



- Molecule 3: THROMBIN HEAVY CHAIN



- Molecule 4: THROMBIN LIGHT CHAIN



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

Chain C:  50% 50%

NAG1
NAG2

- Molecule 6: N-acetyl-alpha-neuraminic acid-(2-6)-beta-D-galactopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-2)-alpha-D-mannopyranose-(1-6)-beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain D:  14% 57% 29%

NAG1
NAG2
BMA3
MAN4
NAG5
GAL6
SIA7

4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	66.77Å 92.23Å 147.49Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	60.83 – 1.90 60.83 – 1.90	Depositor EDS
% Data completeness (in resolution range)	99.6 (60.83-1.90) 99.6 (60.83-1.90)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.07 (at 1.90Å)	Xtriage
Refinement program	REFMAC 5.7.0029	Depositor
R, R_{free}	0.201 , 0.244 0.204 , 0.247	Depositor DCC
R_{free} test set	3744 reflections (5.16%)	wwPDB-VP
Wilson B-factor (Å ²)	30.2	Xtriage
Anisotropy	0.128	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 35.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	6373	wwPDB-VP
Average B, all atoms (Å ²)	35.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.09% 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

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: BMA, SIA, CIT, 0G6, PO4, NAG, PEG, MAN, GAL, NA, PCA

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	0.39	0/1707	0.73	0/2326
2	B	0.42	1/1761 (0.1%)	0.70	1/2415 (0.0%)
3	H	0.46	0/2146	0.73	0/2903
4	L	0.43	0/273	0.73	0/364
All	All	0.43	1/5887 (0.0%)	0.72	1/8008 (0.0%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	1	PCA	C-N	-5.17	1.26	1.33

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	215	TYR	CB-CA-C	-5.10	110.68	116.54

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	1670	0	1583	32	0
2	B	1722	0	1594	46	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	H	2091	0	2045	20	0
4	L	270	0	253	5	0
5	C	28	0	25	1	0
6	D	95	0	80	10	0
7	A	7	0	10	0	0
7	H	14	0	20	0	0
8	B	5	0	0	2	0
8	H	20	0	0	1	0
9	H	13	0	5	4	0
10	H	1	0	0	0	0
11	H	30	0	31	0	0
12	A	95	0	0	2	0
12	B	105	0	0	5	0
12	H	189	0	0	5	1
12	L	18	0	0	0	0
All	All	6373	0	5646	108	1

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 (108) 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)
9:H:608:CIT:O2	9:H:608:CIT:H42	1.54	1.02
1:A:91[A]:ARG:NH1	2:B:111:TYR:CE2	2.37	0.93
12:H:755:HOH:O	4:L:285:THR:HG22	1.71	0.91
6:D:2:NAG:H3	6:D:3:BMA:H2	1.60	0.83
1:A:192:VAL:HG22	1:A:211:LEU:CD2	2.11	0.80
2:B:127:THR:O	2:B:129:VAL:HG23	1.81	0.79
9:H:608:CIT:O2	9:H:608:CIT:C4	2.31	0.78
6:D:5:NAG:H83	6:D:7:SIA:H112	1.65	0.77
6:D:2:NAG:H4	6:D:3:BMA:O2	1.84	0.77
2:B:4:LEU:HD22	2:B:24[B]:VAL:HG22	1.67	0.76
1:A:184:LYS:O	1:A:188:GLU:HG2	1.88	0.73
1:A:109:ARG:HD3	1:A:110:THR:O	1.89	0.72
2:B:3:GLN:CD	2:B:5:ILE:HD11	2.15	0.71
1:A:91[A]:ARG:NH1	2:B:111:TYR:CD2	2.57	0.70
1:A:91[A]:ARG:O	1:A:91[A]:ARG:HG3	1.89	0.70
1:A:164:VAL:HG22	1:A:176:LEU:HD12	1.75	0.68
1:A:106:GLU:HG2	1:A:107:ILE:N	2.06	0.68
2:B:30:THR:OG1	2:B:74:ARG:HD3	1.94	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:84:ASN:OD1	12:B:402:HOH:O	2.13	0.67
2:B:47:TRP:CZ2	2:B:49:GLY:HA2	2.31	0.65
6:D:1:NAG:O4	6:D:2:NAG:C6	2.44	0.65
2:B:73:ASP:O	2:B:76:THR:HG22	1.96	0.64
1:A:192:VAL:HG22	1:A:211:LEU:HD22	1.80	0.63
2:B:4:LEU:CD2	2:B:24[B]:VAL:HG22	2.29	0.63
1:A:167:GLN:HG3	1:A:174:TYR:CZ	2.33	0.63
4:L:306:ASP:CG	4:L:309:GLU:HG2	2.24	0.62
2:B:142:GLN:NE2	12:B:401:HOH:O	2.03	0.62
2:B:132:LYS:N	2:B:132:LYS:HE3	2.14	0.61
2:B:2:VAL:HG21	2:B:27:TYR:CE2	2.35	0.61
1:A:54:ARG:HG2	1:A:58:ILE:HB	1.83	0.60
1:A:91[B]:ARG:HG3	1:A:91[B]:ARG:O	2.01	0.60
9:H:608:CIT:H22	12:H:706:HOH:O	2.02	0.59
2:B:185:SER:O	8:B:301:PO4:O2	2.21	0.59
6:D:2:NAG:C3	6:D:3:BMA:H2	2.29	0.58
2:B:159:GLN:HG2	2:B:160:GLU:HG2	1.86	0.57
6:D:1:NAG:O4	6:D:2:NAG:H62	2.03	0.57
2:B:103:GLU:CD	3:H:388:ARG:HH21	2.14	0.56
2:B:30:THR:HG23	2:B:74:ARG:NH1	2.22	0.55
2:B:207[A]:SER:OG	2:B:223:THR:CG2	2.54	0.55
12:A:401:HOH:O	5:C:1:NAG:O6	2.18	0.55
2:B:146:ASN:HD22	2:B:198:PRO:HA	1.72	0.54
3:H:531:MET:SD	3:H:542:MET:HG3	2.47	0.54
6:D:7:SIA:O1B	6:D:7:SIA:H6	2.07	0.54
1:A:194:ALA:HB2	1:A:209:SER:CB	2.39	0.53
2:B:2:VAL:HG21	2:B:27:TYR:HE2	1.72	0.53
1:A:91[A]:ARG:NH1	2:B:111:TYR:CZ	2.77	0.52
1:A:117:PHE:CZ	2:B:143:PRO:HG3	2.44	0.52
1:A:194:ALA:HB2	1:A:209:SER:HB2	1.92	0.52
2:B:72:GLU:OE2	12:B:404:HOH:O	2.19	0.51
1:A:141:TYR:CG	1:A:142:PRO:HA	2.46	0.51
2:B:186:GLY:HA3	8:B:301:PO4:O1	2.10	0.50
3:H:417:ASP:OD2	8:H:613:PO4:O1	2.28	0.50
1:A:91[A]:ARG:O	1:A:94:TRP:HZ3	1.95	0.50
2:B:3:GLN:OE1	2:B:5:ILE:HD11	2.11	0.50
2:B:85:ASN:O	12:B:403:HOH:O	2.17	0.50
6:D:1:NAG:C6	6:D:2:NAG:H61	2.42	0.50
3:H:474:LYS:HD2	3:H:476:GLN:O	2.12	0.49
1:A:67:SER:HA	1:A:71:PHE:CE2	2.48	0.48
2:B:207[A]:SER:OG	2:B:223:THR:HG22	2.14	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:37:GLN:HB2	1:A:47:LEU:HD11	1.96	0.48
2:B:73:ASP:HB3	2:B:76:THR:HG22	1.95	0.47
1:A:132:SER:HA	1:A:180:LEU:O	2.14	0.47
3:H:363:HIS:O	3:H:372:LYS:HE3	2.14	0.47
2:B:55:ASP:HA	3:H:393:ARG:HD3	1.97	0.47
2:B:13:LYS:O	2:B:14:PRO:C	2.54	0.47
6:D:2:NAG:C3	6:D:3:BMA:C2	2.91	0.46
1:A:117:PHE:HZ	2:B:143:PRO:HG3	1.80	0.46
2:B:142:GLN:OE1	2:B:148:VAL:N	2.43	0.46
4:L:297:PRO:HA	4:L:301:LYS:HB2	1.98	0.46
2:B:81:MET:HE1	2:B:94:TYR:CD2	2.51	0.46
3:H:359:LEU:HD11	3:H:420:ILE:HD11	1.98	0.46
3:H:433:ASP:OD2	4:L:298:LEU:HD11	2.16	0.45
9:H:608:CIT:O3	9:H:608:CIT:O7	2.26	0.45
2:B:97:THR:OG1	2:B:112:PHE:HB3	2.15	0.45
3:H:494:LYS:HE3	12:H:820:HOH:O	2.17	0.45
6:D:3:BMA:O5	6:D:4:MAN:O5	2.34	0.45
2:B:157:PHE:CG	2:B:158:PRO:HA	2.51	0.45
2:B:160:GLU:HB3	2:B:161:PRO:HA	1.99	0.45
3:H:472:VAL:HG12	3:H:473:GLY:N	2.31	0.44
2:B:168:GLU:OE2	2:B:206:LYS:NZ	2.42	0.44
1:A:33:LEU:HD22	1:A:89:GLN:O	2.17	0.44
1:A:54:ARG:HD3	1:A:58:ILE:O	2.18	0.44
2:B:207[A]:SER:OG	2:B:223:THR:HG23	2.17	0.44
1:A:144:GLU:O	1:A:199:HIS:CD2	2.71	0.44
2:B:197:LEU:HD23	2:B:197:LEU:N	2.32	0.43
3:H:363:HIS:CE1	3:H:525:SER:OG	2.71	0.43
1:A:202:LEU:O	1:A:203:SER:C	2.62	0.43
2:B:2:VAL:CG2	2:B:27:TYR:CD2	3.02	0.43
3:H:472:VAL:CG1	3:H:473:GLY:N	2.82	0.43
1:A:109:ARG:CD	1:A:110:THR:O	2.61	0.42
1:A:31[B]:SER:O	1:A:50:ASP:HA	2.18	0.42
2:B:159:GLN:HG3	2:B:180:PRO:HG2	2.02	0.42
3:H:368:PRO:HG2	3:H:412:TRP:CE2	2.54	0.42
3:H:456:GLY:HA3	3:H:529:PHE:CE1	2.54	0.42
1:A:91[A]:ARG:O	1:A:94:TRP:CZ3	2.72	0.42
2:B:103:GLU:CD	3:H:388:ARG:NH2	2.78	0.42
2:B:143:PRO:O	2:B:144:ASP:C	2.62	0.42
1:A:144:GLU:H	1:A:144:GLU:CD	2.27	0.41
2:B:214:HIS:O	2:B:215:TYR:C	2.62	0.41
2:B:73:ASP:CB	2:B:76:THR:HG22	2.50	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:H:357:TRP:CD2	3:H:574:ILE:HG22	2.55	0.41
3:H:439:CYS:SG	4:L:293:CYS:C	3.03	0.41
1:A:144:GLU:O	1:A:199:HIS:HD2	2.04	0.41
3:H:455:LYS:HA	3:H:486:PRO:HA	2.03	0.41
3:H:490:ARG:HD3	12:H:769:HOH:O	2.20	0.40
2:B:5:ILE:HG12	12:B:428:HOH:O	2.21	0.40
1:A:31[B]:SER:OG	12:A:402:HOH:O	2.22	0.40
3:H:399:SER:HB3	12:H:854:HOH:O	2.21	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
12:H:874:HOH:O	12:H:888:HOH:O[4_466]	2.14	0.06

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

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	218/215 (101%)	208 (95%)	10 (5%)	0	100	100
2	B	230/229 (100%)	219 (95%)	11 (5%)	0	100	100
3	H	258/259 (100%)	247 (96%)	11 (4%)	0	100	100
4	L	32/36 (89%)	31 (97%)	1 (3%)	0	100	100
All	All	738/739 (100%)	705 (96%)	33 (4%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	182/187 (97%)	177 (97%)	5 (3%)	39	34
2	B	184/196 (94%)	180 (98%)	4 (2%)	45	42
3	H	223/225 (99%)	222 (100%)	1 (0%)	84	87
4	L	29/31 (94%)	27 (93%)	2 (7%)	14	7
All	All	618/639 (97%)	606 (98%)	12 (2%)	48	47

All (12) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	33	LEU
1	A	106	GLU
1	A	109	ARG
1	A	182	LEU
1	A	190	GLN
2	B	13	LYS
2	B	67	ARG
2	B	132	LYS
2	B	174	THR
3	H	489	GLU
4	L	298	LEU
4	L	301	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	190	GLN
1	A	200	GLN
2	B	146	ASN
3	H	537	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

1 non-standard protein/DNA/RNA residue is 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	$\# Z > 2$
2	PCA	B	1	2	7,8,9	0.59	0	9,10,12	1.43	1 (11%)

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
2	PCA	B	1	2	-	0/0/11/13	0/1/1/1

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	1	PCA	CB-CA-C	-3.03	108.50	112.66

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.5 Carbohydrates ⓘ

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	0.56	0	17,19,21	0.79	0
5	NAG	C	2	5	14,14,15	0.50	0	17,19,21	0.61	0
6	NAG	D	1	6,3	14,14,15	0.49	0	17,19,21	1.71	3 (17%)
6	NAG	D	2	6	14,14,15	0.51	0	17,19,21	0.68	0
6	BMA	D	3	6	11,11,12	0.32	0	15,15,17	0.52	0
6	MAN	D	4	6	11,11,12	0.61	0	15,15,17	0.79	0
6	NAG	D	5	6	14,14,15	0.60	0	17,19,21	1.03	0
6	GAL	D	6	6	11,11,12	0.63	0	15,15,17	0.73	0
6	SIA	D	7	6	20,20,21	0.53	0	21,28,31	1.08	3 (14%)

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	-	0/6/23/26	0/1/1/1
5	NAG	C	2	5	-	2/6/23/26	0/1/1/1
6	NAG	D	1	6,3	-	4/6/23/26	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	MAN	D	4	6	-	2/2/19/22	1/1/1/1
6	NAG	D	5	6	-	0/6/23/26	0/1/1/1
6	GAL	D	6	6	-	1/2/19/22	0/1/1/1
6	SIA	D	7	6	-	4/18/34/38	0/1/1/1

There are no bond length outliers.

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	D	1	NAG	C4-C3-C2	-4.84	103.93	111.02
6	D	1	NAG	C1-O5-C5	2.93	116.11	112.19

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	D	7	SIA	O1B-C1-C2	2.72	119.78	112.71
6	D	1	NAG	O4-C4-C3	2.71	116.76	110.38
6	D	7	SIA	O1A-C1-C2	-2.70	117.01	122.85
6	D	7	SIA	O6-C2-C1	2.20	111.87	107.72

There are no chirality outliers.

All (17) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
6	D	7	SIA	C11-C10-N5-C5
6	D	7	SIA	O10-C10-N5-C5
6	D	1	NAG	O5-C5-C6-O6
6	D	4	MAN	C4-C5-C6-O6
6	D	1	NAG	C4-C5-C6-O6
6	D	3	BMA	O5-C5-C6-O6
6	D	4	MAN	O5-C5-C6-O6
6	D	2	NAG	O5-C5-C6-O6
6	D	7	SIA	C7-C8-C9-O9
6	D	3	BMA	C4-C5-C6-O6
6	D	7	SIA	O8-C8-C9-O9
6	D	1	NAG	C1-C2-N2-C7
6	D	2	NAG	C4-C5-C6-O6
5	C	2	NAG	C3-C2-N2-C7
5	C	2	NAG	C1-C2-N2-C7
6	D	1	NAG	C3-C2-N2-C7
6	D	6	GAL	C4-C5-C6-O6

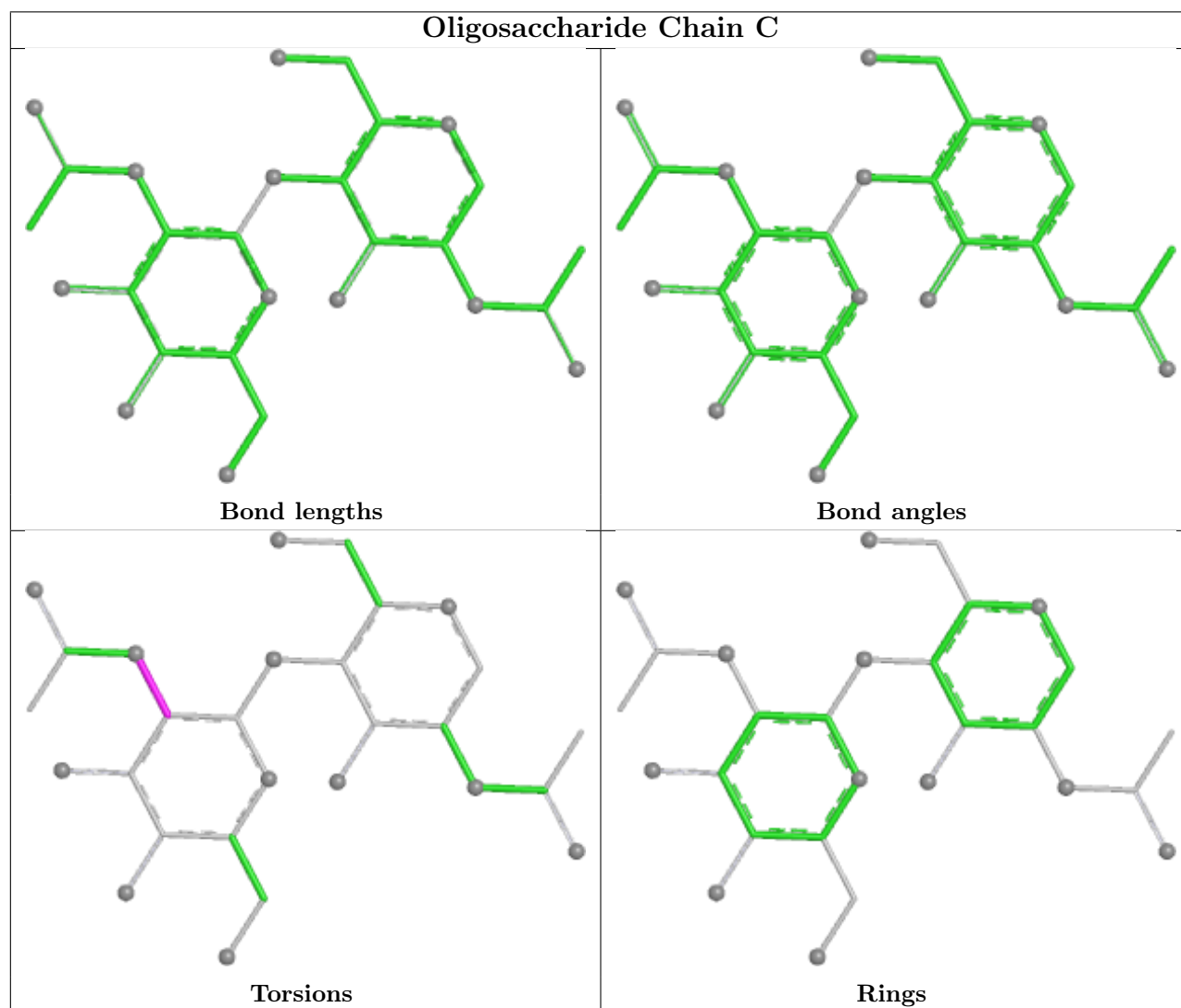
All (1) ring outliers are listed below:

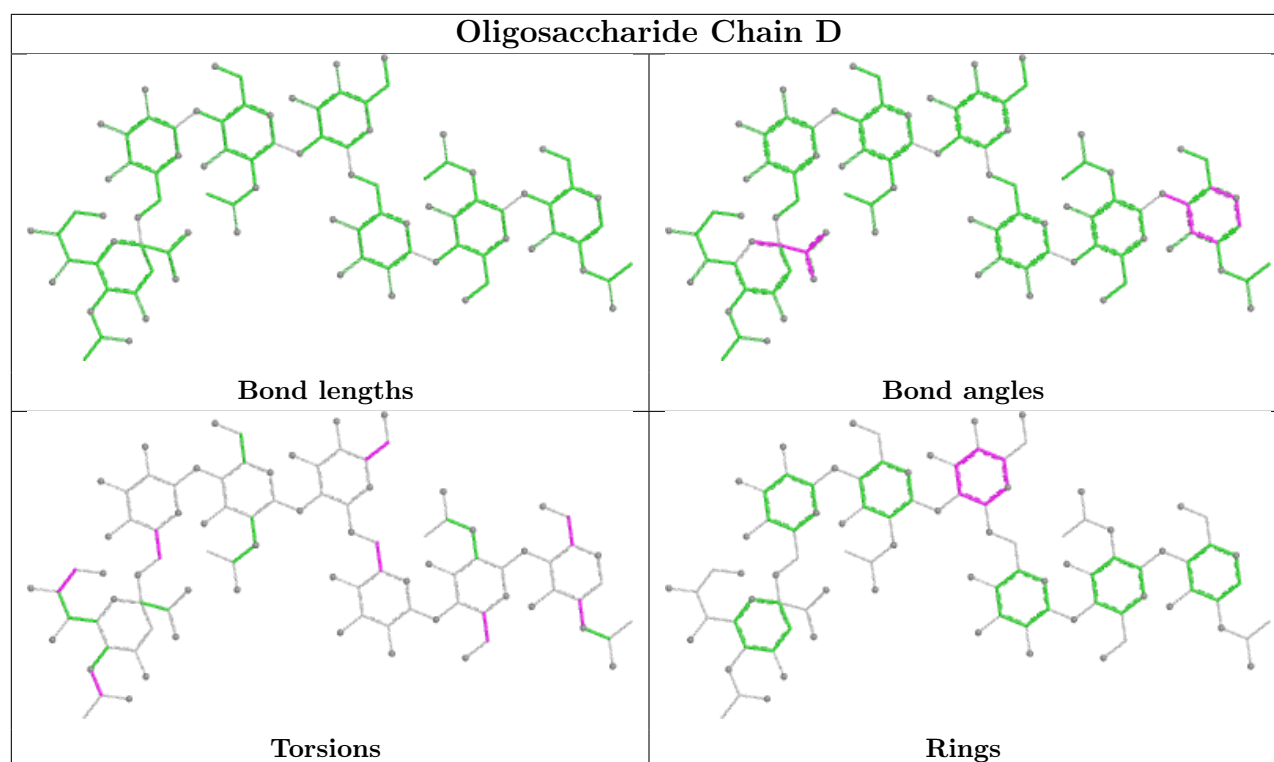
Mol	Chain	Res	Type	Atoms
6	D	4	MAN	C1-C2-C3-C4-C5-O5

7 monomers are involved in 11 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	D	7	SIA	2	0
6	D	4	MAN	1	0
6	D	3	BMA	5	0
6	D	1	NAG	3	0
6	D	2	NAG	7	0
6	D	5	NAG	1	0
5	C	1	NAG	1	0

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





5.6 Ligand geometry [i](#)

Of 11 ligands modelled in this entry, 1 is monoatomic - leaving 10 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	$\# Z > 2$
8	PO4	H	614	-	4,4,4	0.92	0	6,6,6	0.46	0
8	PO4	B	301	-	4,4,4	0.91	0	6,6,6	0.49	0
11	OG6	H	615	-	30,31,32	0.49	0	37,41,42	0.87	2 (5%)
7	PEG	H	609	-	6,6,6	0.51	0	5,5,5	0.28	0
8	PO4	H	611	-	4,4,4	0.83	0	6,6,6	0.56	0
8	PO4	H	612	-	4,4,4	0.73	0	6,6,6	0.48	0
7	PEG	H	616	-	6,6,6	0.42	0	5,5,5	0.37	0
7	PEG	A	303	-	6,6,6	0.46	0	5,5,5	0.33	0
9	CIT	H	608	-	12,12,12	1.17	0	17,17,17	1.21	2 (11%)
8	PO4	H	613	-	4,4,4	0.96	0	6,6,6	0.44	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
11	0G6	H	615	-	-	1/31/41/43	0/2/2/2
7	PEG	H	609	-	-	1/4/4/4	-
7	PEG	H	616	-	-	3/4/4/4	-
7	PEG	A	303	-	-	3/4/4/4	-
9	CIT	H	608	-	-	9/16/16/16	-

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
11	H	615	0G6	O1-C1-N2	-3.09	117.42	122.96
11	H	615	0G6	CA1-C1-N2	2.63	122.30	116.52
9	H	608	CIT	O6-C6-C3	2.41	117.77	113.14
9	H	608	CIT	C3-C4-C5	-2.16	108.00	113.92

There are no chirality outliers.

All (17) torsion outliers are listed below:

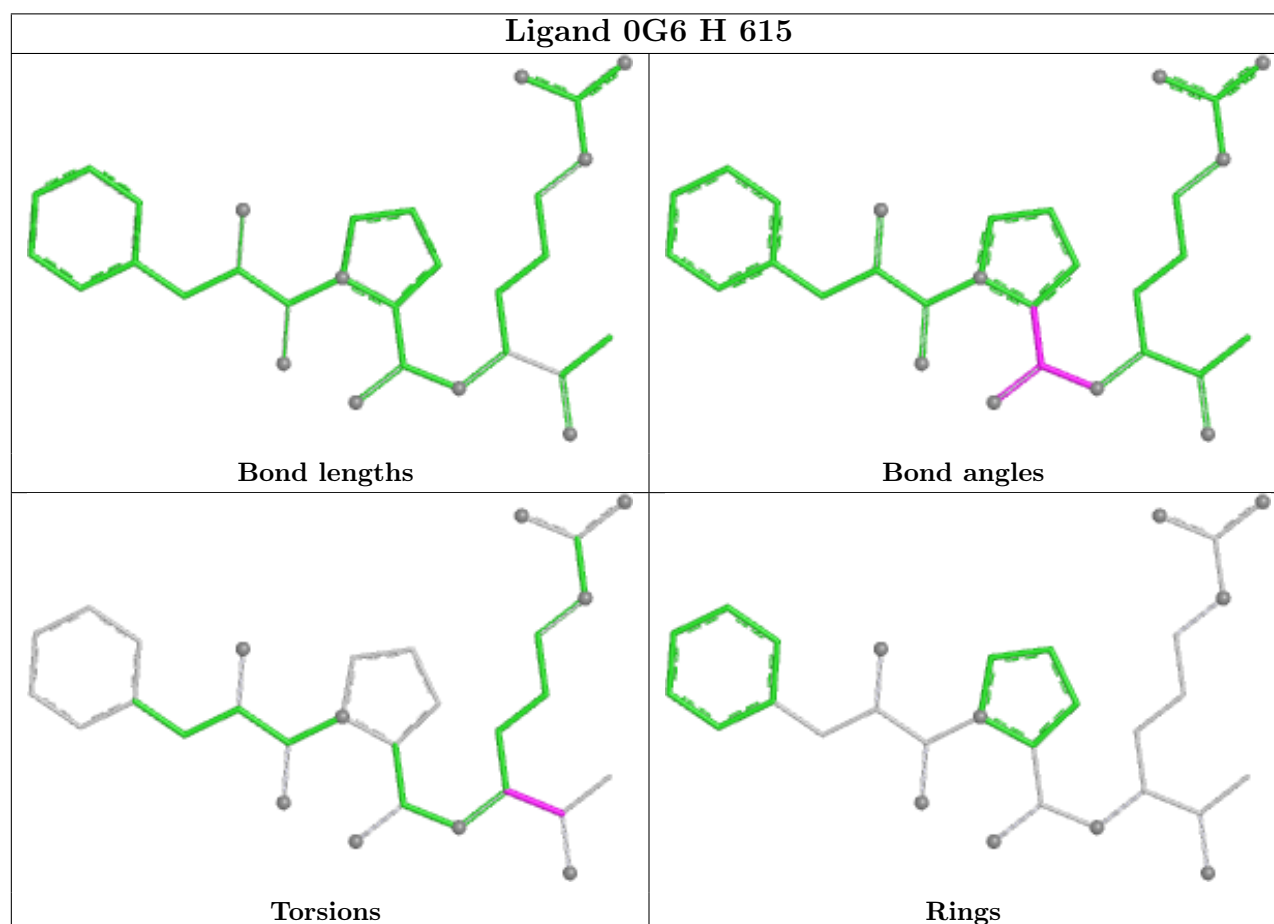
Mol	Chain	Res	Type	Atoms
9	H	608	CIT	C1-C2-C3-O7
9	H	608	CIT	C1-C2-C3-C4
9	H	608	CIT	C2-C3-C6-O5
9	H	608	CIT	C2-C3-C6-O6
9	H	608	CIT	O7-C3-C6-O5
9	H	608	CIT	O7-C3-C6-O6
9	H	608	CIT	C1-C2-C3-C6
7	A	303	PEG	O2-C3-C4-O4
7	A	303	PEG	C1-C2-O2-C3
7	H	616	PEG	C4-C3-O2-C2
7	H	616	PEG	O1-C1-C2-O2
7	A	303	PEG	O1-C1-C2-O2
7	H	609	PEG	C1-C2-O2-C3
9	H	608	CIT	O2-C1-C2-C3
9	H	608	CIT	O1-C1-C2-C3
7	H	616	PEG	C1-C2-O2-C3
11	H	615	0G6	O2-C2-CA2-CB2

There are no ring outliers.

3 monomers are involved in 7 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
8	B	301	PO4	2	0
9	H	608	CIT	4	0
8	H	613	PO4	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 all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.



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	215/215 (100%)	0.21	3 (1%) 73 76	15, 35, 51, 63	5 (2%)
2	B	228/229 (99%)	0.35	8 (3%) 47 50	16, 37, 56, 69	3 (1%)
3	H	259/259 (100%)	-0.30	2 (0%) 82 85	15, 27, 39, 51	1 (0%)
4	L	33/36 (91%)	0.17	1 (3%) 52 56	18, 33, 47, 54	1 (3%)
All	All	735/739 (99%)	0.07	14 (1%) 66 70	15, 33, 51, 69	10 (1%)

All (14) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	141	THR	4.2
2	B	229	PRO	2.8
3	H	472	VAL	2.8
1	A	154	ALA	2.7
2	B	185	SER	2.7
2	B	140	SER	2.6
2	B	228	VAL	2.5
4	L	317	ILE	2.4
1	A	155	LEU	2.4
3	H	515	GLY	2.3
2	B	75	SER	2.3
1	A	204	SER	2.2
2	B	197	LEU	2.2
2	B	73	ASP	2.1

6.2 Non-standard residues in protein, DNA, RNA chains [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($\text{\AA}^2$)	Q<0.9
2	PCA	B	1	8/9	0.82	0.14	54,55,58,59	0

6.3 Carbohydrates

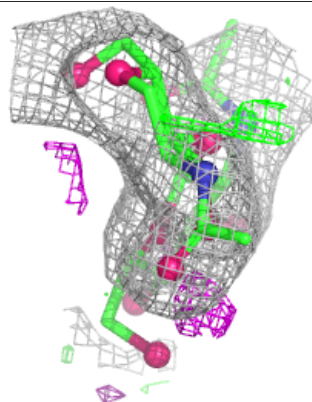
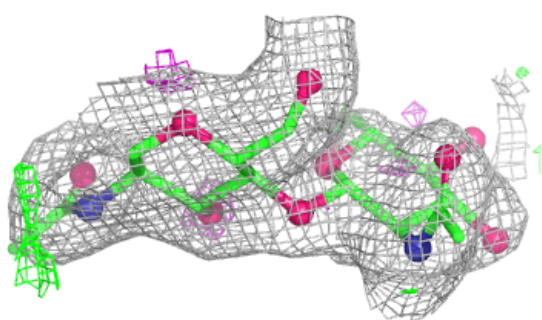
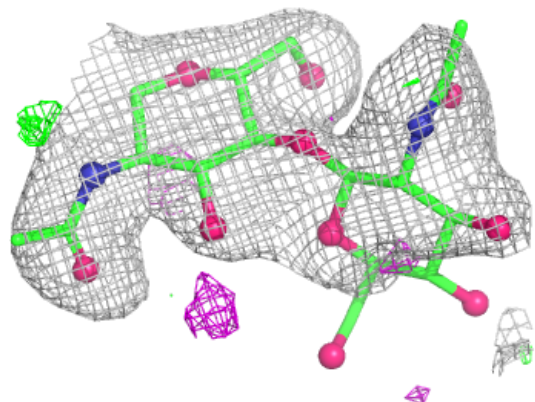
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
5	NAG	C	1	14/15	-	-	46,50,56,66	0
5	NAG	C	2	14/15	-	-	75,80,84,87	0
6	NAG	D	1	14/15	-	-	54,61,66,76	0
6	NAG	D	2	14/15	-	-	86,91,96,103	0
6	BMA	D	3	11/12	-	-	90,103,106,109	0
6	MAN	D	4	11/12	-	-	80,85,87,88	0
6	NAG	D	5	14/15	-	-	50,63,74,75	0
6	GAL	D	6	11/12	-	-	43,44,45,46	0
6	SIA	D	7	20/21	-	-	47,61,68,69	0

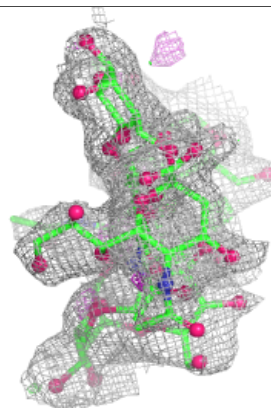
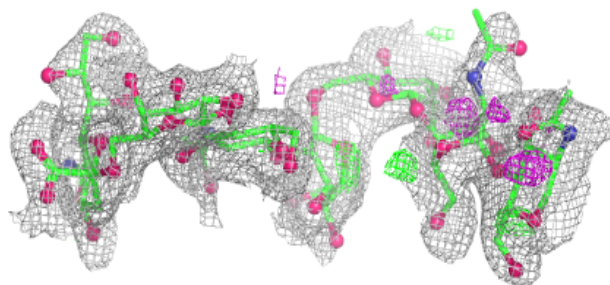
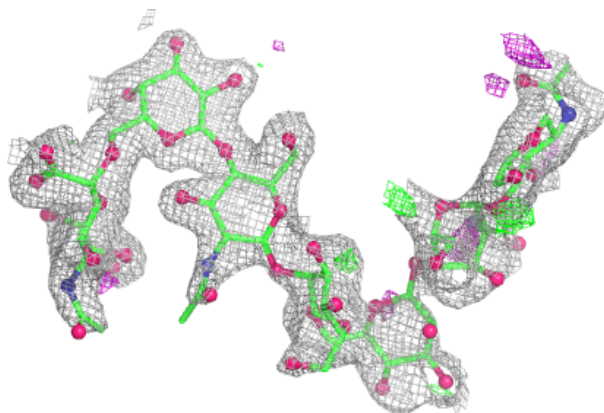
The following is a graphical depiction of the model fit to experimental electron density for oligosaccharide. Each fit is shown from different orientation to approximate a three-dimensional view.

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)



6.4 Ligands ⓘ

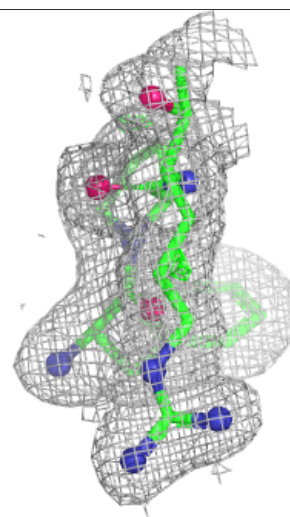
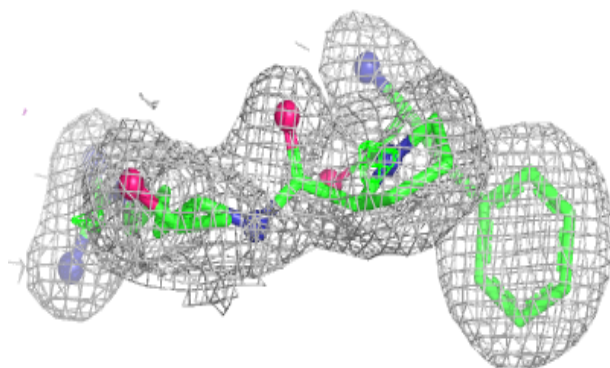
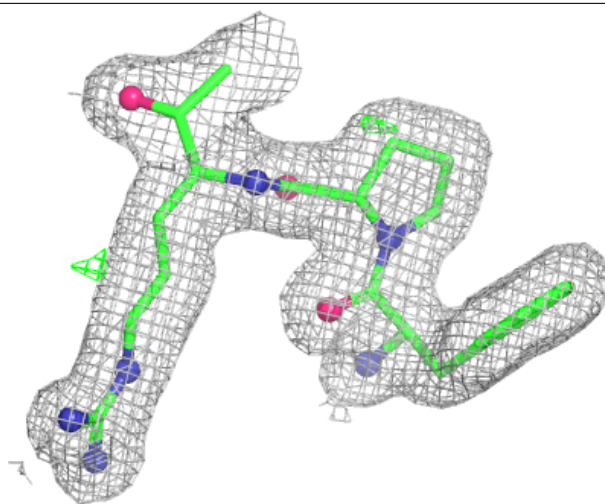
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q<0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
8	PO4	H	614	5/5	0.71	0.14	94,96,99,100	0
8	PO4	B	301	5/5	0.74	0.15	100,104,105,106	0
9	CIT	H	608	13/13	0.78	0.15	50,55,63,66	0
8	PO4	H	613	5/5	0.83	0.11	77,79,81,83	0
7	PEG	H	609	7/7	0.84	0.15	34,43,47,51	0
7	PEG	H	616	7/7	0.87	0.12	52,53,55,56	0
7	PEG	A	303	7/7	0.88	0.13	50,52,53,53	0
8	PO4	H	611	5/5	0.94	0.08	53,54,57,58	0
10	NA	H	610	1/1	0.97	0.06	34,34,34,34	0
11	OG6	H	615	30/31	0.97	0.05	19,21,22,23	0
8	PO4	H	612	5/5	0.98	0.10	32,33,33,34	0

The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

Electron density around 0G6 H 615:

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



6.5 Other polymers [i](#)

There are no such residues in this entry.