



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 18, 2026 – 10:05 AM UTC

PDB ID : 6LXI / pdb_00006lxi
Title : Crystal structure of Z2B3 Fab in complex with influenza virus neuraminidase from A/Brevig Mission/1/1918 (H1N1)
Authors : Jiang, H.; Peng, W.; Qi, J.; Chai, Y.; Song, H.; Shi, Y.; Gao, G.F.; Wu, Y.
Deposited on : 2020-02-11
Resolution : 2.50 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:

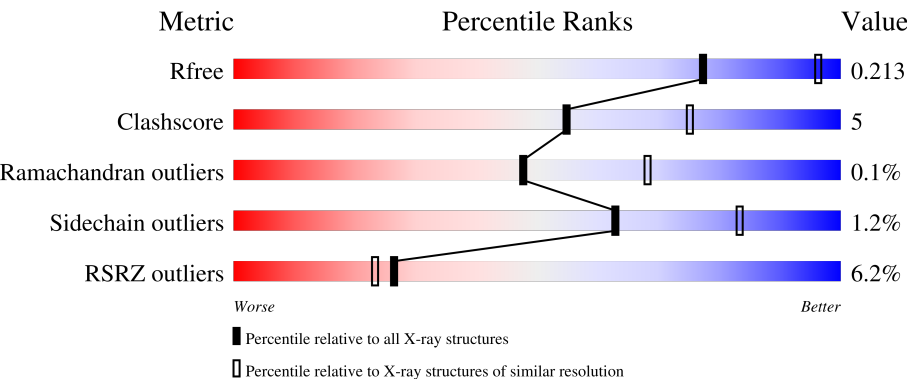
MolProbity : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 2.0
EDS : 3.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (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 2.50 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	5829 (2.50-2.50)
Clashscore	190562	6492 (2.50-2.50)
Ramachandran outliers	187476	6378 (2.50-2.50)
Sidechain outliers	187428	6380 (2.50-2.50)
RSRZ outliers	180081	5833 (2.50-2.50)

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	469	78%17%
1	B	469	75%8%17%
2	C	237	8% 78%12%10%
2	H	237	% 92%5%
3	D	216	25% 69%20%9%

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Mol	Chain	Length	Quality of chain
3	L	216	12% 80% 18%
4	E	5	20% 60% 20%
4	F	5	60% 40%

2 Entry composition [i](#)

There are 6 unique types of molecules in this entry. The entry contains 12997 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 Neuraminidase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	387	Total	C	N	O	S	0	0	0
			2980	1866	512	579	23			
1	B	387	Total	C	N	O	S	0	0	0
			2980	1866	512	579	23			

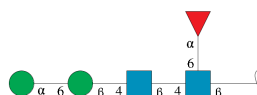
- Molecule 2 is a protein called Heavy chain of Z2B3 Fab.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	H	232	Total	C	N	O	S	0	0	0
			1727	1091	281	348	7			
2	C	214	Total	C	N	O	S	0	0	0
			1611	1023	261	320	7			

- Molecule 3 is a protein called Light chain of Z2B3 Fab.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	L	211	Total	C	N	O	S	0	0	0
			1563	973	260	325	5			
3	D	196	Total	C	N	O	S	0	0	0
			1447	902	237	303	5			

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



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
4	E	5	Total	C	N	O	0	0	0
			60	34	2	24			

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
4	F	5	Total	C	N	O	0	0	0
			60	34	2	24			

- Molecule 5 is CALCIUM ION (CCD ID: CA) (formula: Ca) (labeled as "Ligand of Interest" by depositor).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	2	Total	Ca	0	0
			2	2		
5	B	2	Total	Ca	0	0
			2	2		

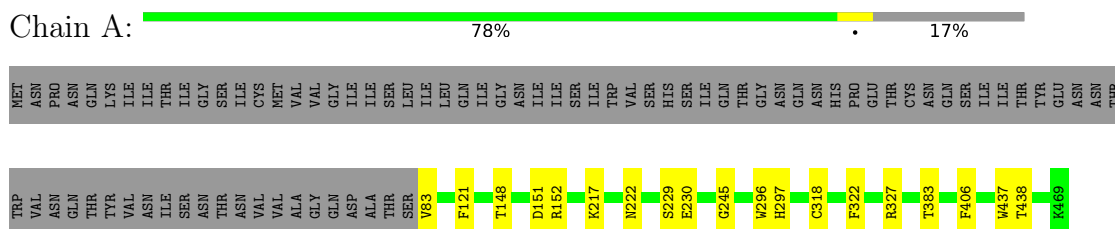
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	221	Total	O	0	0
			221	221		
6	H	69	Total	O	0	0
			69	69		
6	L	46	Total	O	0	0
			46	46		
6	B	165	Total	O	0	0
			165	165		
6	C	46	Total	O	0	0
			46	46		
6	D	18	Total	O	0	0
			18	18		

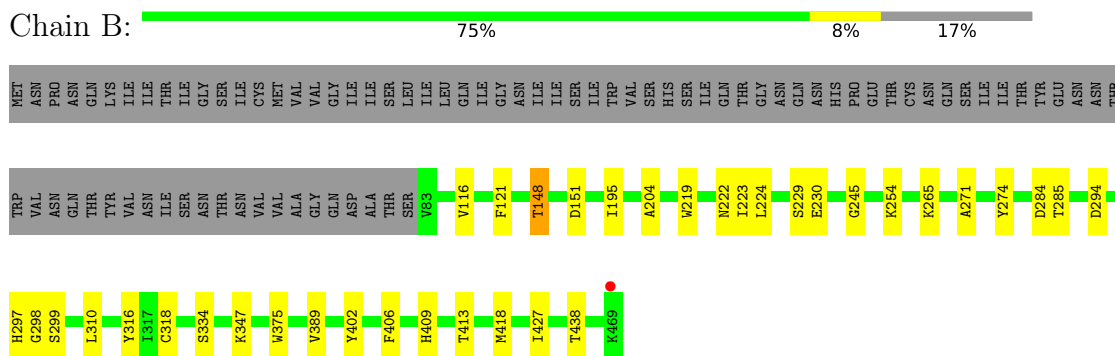
3 Residue-property plots

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.

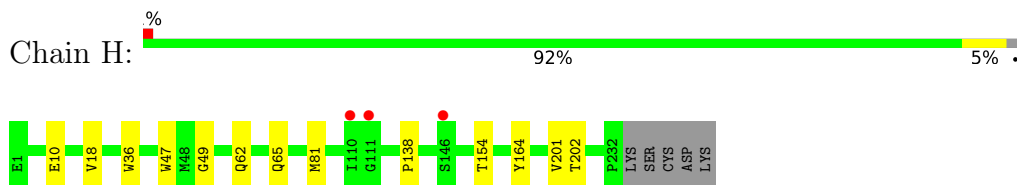
- Molecule 1: Neuraminidase



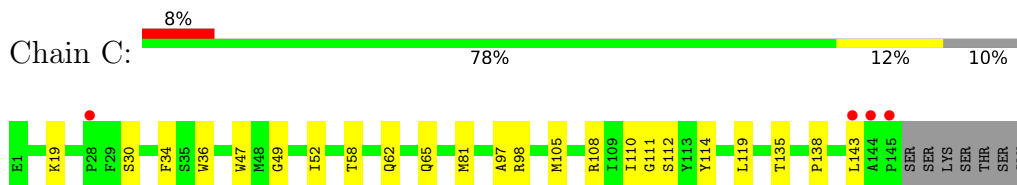
- Molecule 1: Neuraminidase

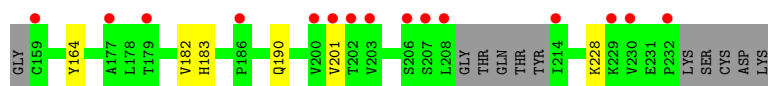


- Molecule 2: Heavy chain of Z2B3 Fab

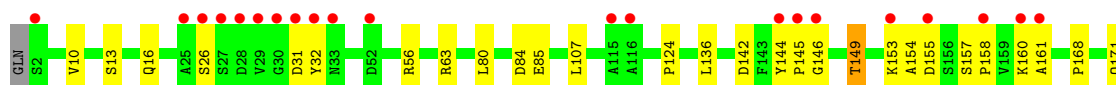
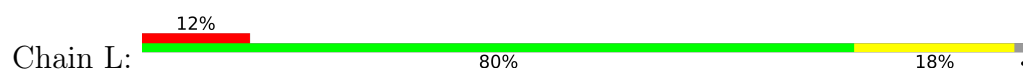


- Molecule 2: Heavy chain of Z2B3 Fab





- Molecule 3: Light chain of Z2B3 Fab



- Molecule 3: Light chain of Z2B3 Fab



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



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



4 Data and refinement statistics

Property	Value	Source
Space group	P 4 21 2	Depositor
Cell constants a, b, c, α , β , γ	163.13Å 163.13Å 190.76Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	47.93 – 2.50 47.93 – 2.50	Depositor EDS
% Data completeness (in resolution range)	99.0 (47.93-2.50) 99.0 (47.93-2.50)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.06 (at 2.51Å)	Xtriage
Refinement program	PHENIX 1.18.2-3874_1692	Depositor
R, R_{free}	0.183 , 0.212 0.186 , 0.213	Depositor DCC
R_{free} test set	4445 reflections (4.97%)	wwPDB-VP
Wilson B-factor (Å ²)	38.1	Xtriage
Anisotropy	0.077	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 39.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	12997	wwPDB-VP
Average B, all atoms (Å ²)	41.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.54% 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: BMA, NAG, CA, MAN, FUC

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	0.65	2/3061 (0.1%)	0.62	0/4160
1	B	0.64	3/3061 (0.1%)	0.62	0/4160
2	C	0.46	0/1650	0.51	0/2249
2	H	0.50	0/1769	0.49	0/2413
3	D	0.54	2/1479 (0.1%)	0.70	1/2018 (0.0%)
3	L	0.49	0/1600	0.59	0/2183
All	All	0.57	7/12620 (0.1%)	0.60	1/17183 (0.0%)

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	297	HIS	C-O	-7.73	1.17	1.24
3	D	54	SER	C-N	-7.11	1.23	1.33
1	A	296	TRP	C-O	-6.51	1.18	1.23
1	B	297	HIS	C-O	-6.25	1.18	1.24
1	B	298	GLY	C-N	-5.52	1.25	1.33
1	B	224	LEU	C-O	-5.46	1.17	1.23
3	D	52	ASP	C-O	5.17	1.31	1.23

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	D	176	TYR	N-CA-C	5.99	119.58	109.76

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	2980	0	2802	10	0
1	B	2980	0	2802	21	0
2	C	1611	0	1573	18	0
2	H	1727	0	1686	7	0
3	D	1447	0	1385	34	0
3	L	1563	0	1505	24	0
4	E	60	0	52	4	0
4	F	60	0	52	3	0
5	A	2	0	0	0	0
5	B	2	0	0	0	0
6	A	221	0	0	1	1
6	B	165	0	0	2	1
6	C	46	0	0	0	0
6	D	18	0	0	0	1
6	H	69	0	0	0	0
6	L	46	0	0	2	0
All	All	12997	0	11857	116	2

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 5.

All (116) 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)
4:F:3:BMA:H62	4:F:4:MAN:H5	1.55	0.85
3:D:135:THR:HG22	3:D:183:SER:HA	1.69	0.75
4:E:3:BMA:H62	4:E:4:MAN:H5	1.68	0.74
3:L:171:GLN:OE1	3:L:173:ASN:ND2	2.21	0.73
1:B:413:THR:O	6:B:601:HOH:O	2.05	0.73
3:L:171:GLN:HE21	3:L:177:ALA:HB2	1.53	0.72
3:D:127:GLU:HG2	3:D:130:GLN:NE2	2.07	0.70
1:B:230:GLU:OE1	6:B:602:HOH:O	2.11	0.69
3:D:127:GLU:O	3:D:130:GLN:NE2	2.24	0.68
3:D:130:GLN:N	3:D:130:GLN:OE1	2.26	0.68
4:E:3:BMA:C6	4:E:4:MAN:H5	2.23	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:L:173:ASN:OD1	3:L:175:LYS:HG3	1.93	0.68
3:D:136:LEU:HD12	3:D:182:LEU:HD23	1.77	0.67
2:C:182:VAL:HG12	2:C:201:VAL:HG22	1.78	0.66
3:L:171:GLN:HB2	3:L:173:ASN:ND2	2.12	0.64
3:L:136:LEU:HD12	3:L:182:LEU:HD23	1.79	0.64
2:C:110:ILE:HG22	2:C:111:GLY:H	1.65	0.62
3:D:200:THR:HB	3:D:205:THR:HG23	1.81	0.61
1:B:148:THR:CG2	1:B:438:THR:H	2.13	0.61
3:D:170:LYS:HD2	3:D:174:ASN:OD1	1.99	0.61
3:D:185:THR:N	3:D:188:GLN:OE1	2.33	0.60
2:C:143:LEU:HD13	3:D:137:VAL:HG21	1.81	0.60
3:D:144:TYR:CD1	3:D:145:PRO:HA	2.37	0.59
1:A:148:THR:CG2	1:A:438:THR:H	2.15	0.59
3:D:149:THR:HB	3:D:200:THR:CG2	2.33	0.58
1:A:217:LYS:O	6:A:603:HOH:O	2.17	0.58
2:C:36:TRP:CE2	2:C:81:MET:HB2	2.39	0.58
2:C:138:PRO:HB3	2:C:164:TYR:HB3	1.85	0.58
3:L:144:TYR:HD1	3:L:145:PRO:CA	2.17	0.57
3:L:160:LYS:HG3	3:L:161:ALA:N	2.18	0.57
3:L:153:LYS:HG2	3:L:158:PRO:HA	1.85	0.57
3:L:155:ASP:OD1	3:L:193:ARG:HB3	2.04	0.57
1:B:223:ILE:O	1:B:223:ILE:HG22	2.05	0.57
1:B:284:ASP:OD1	1:B:285:THR:HG23	2.04	0.56
1:B:427:ILE:HG21	2:C:105:MET:HE1	1.86	0.56
1:B:148:THR:HG22	1:B:438:THR:H	1.71	0.55
1:B:222:ASN:HB3	1:B:245:GLY:HA2	1.88	0.55
1:A:230:GLU:OE2	1:A:406:PHE:HA	2.07	0.55
3:L:26:SER:O	3:L:31:ASP:HB2	2.06	0.55
2:H:36:TRP:CE2	2:H:81:MET:HB2	2.41	0.55
2:C:58:THR:O	2:C:112:SER:HA	2.06	0.55
2:C:62:GLN:HA	2:C:65:GLN:HG3	1.89	0.55
2:C:110:ILE:HG22	2:C:111:GLY:N	2.24	0.53
1:B:195:ILE:HG12	1:B:204:ALA:HB2	1.89	0.53
1:A:148:THR:HG22	1:A:437:TRP:HA	1.90	0.53
2:H:10:GLU:HG2	2:H:18:VAL:HG23	1.91	0.53
1:B:265:LYS:HG2	1:B:310:LEU:HD12	1.92	0.52
2:H:62:GLN:HA	2:H:65:GLN:HG3	1.92	0.52
1:B:294:ASP:OD2	1:B:316:TYR:OH	2.27	0.51
3:D:117:PRO:HD3	3:D:201:HIS:ND1	2.24	0.51
3:L:63:ARG:NH2	3:L:84:ASP:OD2	2.44	0.51
3:L:154:ALA:HB2	3:L:195:TYR:CE2	2.46	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:D:146:GLY:O	3:D:168:PRO:HG2	2.11	0.51
3:L:31:ASP:OD2	6:L:302:HOH:O	2.19	0.51
2:C:19:LYS:NZ	4:F:4:MAN:O5	2.44	0.50
3:D:140:ILE:HD11	3:D:150:VAL:HG21	1.93	0.50
3:D:201:HIS:O	3:D:203:GLY:O	2.29	0.49
2:C:228:LYS:HE2	2:C:228:LYS:HA	1.94	0.49
4:E:1:NAG:H62	4:E:2:NAG:C7	2.43	0.49
3:D:112:GLN:NE2	3:D:144:TYR:CD2	2.81	0.49
1:A:151:ASP:OD2	1:A:152:ARG:HG3	2.13	0.48
3:L:144:TYR:HD1	3:L:145:PRO:HA	1.78	0.48
3:D:112:GLN:NE2	3:D:144:TYR:HD2	2.11	0.48
1:B:121:PHE:CG	1:B:229:SER:HA	2.49	0.48
3:D:122:PHE:HB2	3:D:137:VAL:HG22	1.95	0.48
3:D:28:ASP:OD1	3:D:29:VAL:N	2.41	0.47
1:B:409:HIS:CD2	1:B:418:MET:HE1	2.49	0.47
4:F:3:BMA:H61	4:F:4:MAN:H2	1.35	0.47
2:C:47:TRP:CH2	2:C:49:GLY:HA2	2.50	0.47
4:E:3:BMA:H61	4:E:4:MAN:H2	1.42	0.47
3:D:199:VAL:O	3:D:205:THR:HA	2.14	0.47
1:B:116:VAL:HG11	1:B:148:THR:HG21	1.96	0.46
1:B:271:ALA:HB1	1:B:274:TYR:HB2	1.96	0.46
3:L:10:VAL:HG23	3:L:107:LEU:HD13	1.98	0.46
2:H:138:PRO:HB3	2:H:164:TYR:HB3	1.98	0.45
3:D:149:THR:HB	3:D:200:THR:HG23	1.99	0.45
2:C:97:ALA:HB1	2:C:119:LEU:HB3	1.99	0.45
2:H:36:TRP:CD2	2:H:81:MET:HB2	2.52	0.45
3:L:149:THR:HG23	3:L:200:THR:OG1	2.17	0.45
3:D:144:TYR:CD1	3:D:145:PRO:CA	2.99	0.45
3:D:200:THR:HA	3:D:205:THR:HA	1.98	0.45
1:B:347:LYS:HB3	1:B:402:TYR:CG	2.52	0.45
3:L:146:GLY:O	3:L:168:PRO:HG2	2.17	0.44
1:B:219:TRP:CE2	1:B:254:LYS:HE3	2.53	0.44
1:B:151:ASP:OD2	2:C:108:ARG:HD3	2.18	0.44
2:C:183:HIS:CE1	3:D:177:ALA:HB3	2.52	0.44
3:L:31:ASP:OD2	3:L:32:TYR:CE1	2.70	0.44
3:L:124:PRO:HD3	3:L:136:LEU:HD23	1.99	0.43
2:C:52:ILE:HD11	2:C:114:TYR:HA	1.99	0.43
3:L:16:GLN:HB3	6:L:329:HOH:O	2.17	0.43
2:C:36:TRP:CD2	2:C:81:MET:HB2	2.52	0.43
3:D:127:GLU:HG2	3:D:130:GLN:HE22	1.83	0.43
3:D:142:ASP:H	3:D:171:GLN:HE22	1.66	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:D:142:ASP:H	3:D:171:GLN:NE2	2.16	0.43
1:A:148:THR:HG23	1:A:438:THR:H	1.83	0.43
3:D:121:LEU:HD22	3:D:197:CYS:HB2	1.99	0.43
3:D:142:ASP:CA	3:D:175:LYS:HD3	2.49	0.43
1:A:322:PHE:HB2	1:A:327:ARG:HD2	2.00	0.42
3:L:186:PRO:O	3:L:190:LYS:HG2	2.19	0.42
2:C:34:PHE:CZ	2:C:98:ARG:HD2	2.54	0.42
3:D:112:GLN:HG2	3:D:144:TYR:CE2	2.54	0.42
2:H:47:TRP:CH2	2:H:49:GLY:HA2	2.54	0.42
1:A:121:PHE:CG	1:A:229:SER:HA	2.54	0.42
3:D:27:SER:HB2	3:D:95:ARG:HG3	2.01	0.42
2:H:154:THR:OG1	2:H:202:THR:HG22	2.20	0.42
1:B:318:CYS:HA	1:B:334:SER:O	2.20	0.42
1:A:222:ASN:HB3	1:A:245:GLY:HA2	2.01	0.42
3:D:52:ASP:O	3:D:53:VAL:HB	2.19	0.42
3:L:142:ASP:H	3:L:171:GLN:HE22	1.67	0.42
3:D:210:VAL:HG12	3:D:211:ALA:H	1.85	0.41
3:L:85:GLU:HG3	3:L:107:LEU:O	2.21	0.41
1:B:375:TRP:HB3	1:B:389:VAL:HB	2.02	0.41
3:D:202:GLU:HB3	3:D:203:GLY:H	1.75	0.41
1:A:318:CYS:O	1:A:383:THR:HA	2.21	0.40
3:L:207:GLU:OE1	3:L:207:GLU:N	2.54	0.40
1:B:230:GLU:OE2	1:B:406:PHE:HA	2.21	0.40

All (2) 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)
6:B:709:HOH:O	6:D:318:HOH:O[7_555]	2.06	0.14
6:A:719:HOH:O	6:A:739:HOH:O[3_455]	2.13	0.07

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	385/469 (82%)	369 (96%)	16 (4%)	0	100	100
1	B	385/469 (82%)	368 (96%)	17 (4%)	0	100	100
2	C	208/237 (88%)	203 (98%)	5 (2%)	0	100	100
2	H	230/237 (97%)	225 (98%)	5 (2%)	0	100	100
3	D	190/216 (88%)	184 (97%)	5 (3%)	1 (0%)	24	43
3	L	209/216 (97%)	204 (98%)	5 (2%)	0	100	100
All	All	1607/1844 (87%)	1553 (97%)	53 (3%)	1 (0%)	48	68

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
3	D	202	GLU

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	332/407 (82%)	331 (100%)	1 (0%)	86	94
1	B	332/407 (82%)	330 (99%)	2 (1%)	78	91
2	C	183/200 (92%)	180 (98%)	3 (2%)	55	79
2	H	195/200 (98%)	194 (100%)	1 (0%)	81	92
3	D	165/183 (90%)	162 (98%)	3 (2%)	51	77
3	L	178/183 (97%)	172 (97%)	6 (3%)	32	60
All	All	1385/1580 (88%)	1369 (99%)	16 (1%)	63	83

All (16) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	83	VAL
2	H	201	VAL

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Mol	Chain	Res	Type
3	L	13	SER
3	L	56	ARG
3	L	80	LEU
3	L	149	THR
3	L	157	SER
3	L	194	SER
1	B	148	THR
1	B	299	SER
2	C	30	SER
2	C	135	THR
2	C	190	GLN
3	D	76	THR
3	D	80	LEU
3	D	206	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
2	H	117	ASN
2	H	190	GLN
3	L	39	GLN
1	B	409	HIS
2	C	117	ASN
2	C	183	HIS
3	D	39	GLN
3	D	112	GLN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

10 monosaccharides are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
4	NAG	E	1	4,1	14,14,15	0.30	0	17,19,21	0.59	0
4	NAG	E	2	4	14,14,15	0.31	0	17,19,21	0.87	1 (5%)
4	BMA	E	3	4	11,11,12	0.25	0	15,15,17	0.55	0
4	MAN	E	4	4	11,11,12	0.26	0	15,15,17	0.55	0
4	FUC	E	5	4	10,10,11	0.30	0	14,14,16	0.59	0
4	NAG	F	1	4,1	14,14,15	0.32	0	17,19,21	0.56	0
4	NAG	F	2	4	14,14,15	0.30	0	17,19,21	0.61	0
4	BMA	F	3	4	11,11,12	0.27	0	15,15,17	0.52	0
4	MAN	F	4	4	11,11,12	0.26	0	15,15,17	0.57	0
4	FUC	F	5	4	10,10,11	0.30	0	14,14,16	0.58	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
4	NAG	E	1	4,1	-	0/6/23/26	0/1/1/1
4	NAG	E	2	4	-	0/6/23/26	0/1/1/1
4	BMA	E	3	4	-	0/2/19/22	0/1/1/1
4	MAN	E	4	4	-	1/2/19/22	0/1/1/1
4	FUC	E	5	4	-	-	0/1/1/1
4	NAG	F	1	4,1	-	0/6/23/26	0/1/1/1
4	NAG	F	2	4	-	0/6/23/26	0/1/1/1
4	BMA	F	3	4	-	2/2/19/22	0/1/1/1
4	MAN	F	4	4	-	2/2/19/22	0/1/1/1
4	FUC	F	5	4	-	-	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(°)
4	E	2	NAG	O5-C1-C2	-2.44	107.52	111.29

There are no chirality outliers.

All (5) torsion outliers are listed below:

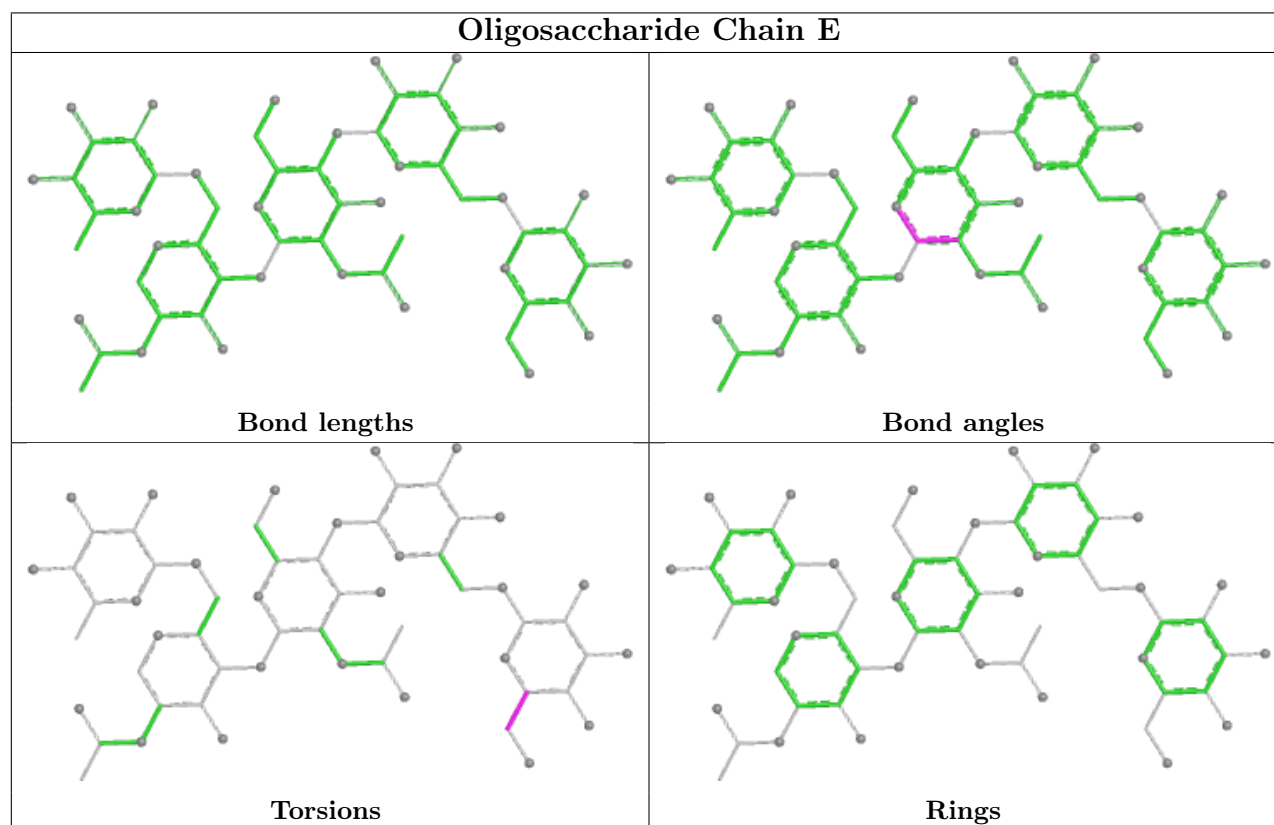
Mol	Chain	Res	Type	Atoms
4	F	3	BMA	C4-C5-C6-O6
4	F	3	BMA	O5-C5-C6-O6
4	F	4	MAN	C4-C5-C6-O6
4	F	4	MAN	O5-C5-C6-O6
4	E	4	MAN	O5-C5-C6-O6

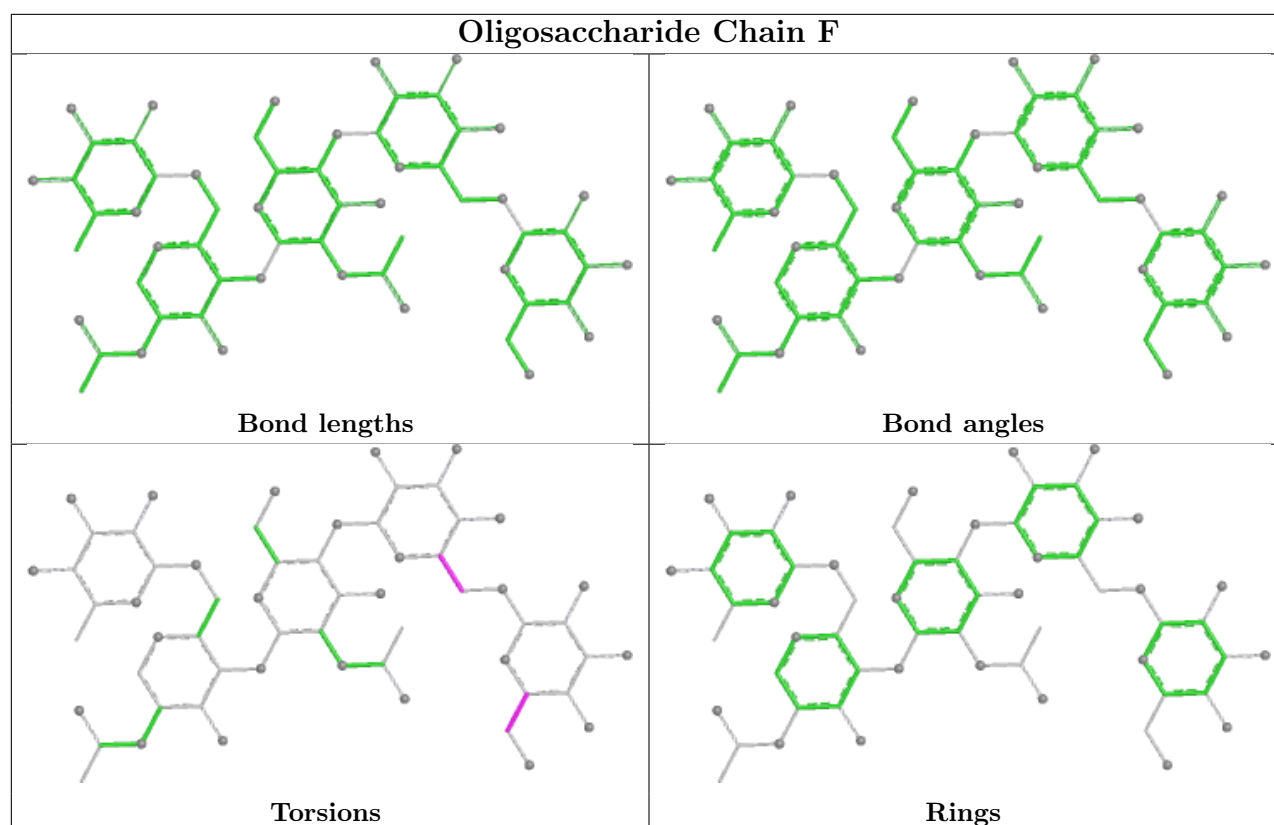
There are no ring outliers.

6 monomers are involved in 7 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	E	3	BMA	3	0
4	F	3	BMA	2	0
4	E	1	NAG	1	0
4	E	4	MAN	3	0
4	E	2	NAG	1	0
4	F	4	MAN	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 4 ligands modelled in this entry, 4 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 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	387/469 (82%)	-0.72	0 100 100	17, 26, 35, 60	0
1	B	387/469 (82%)	-0.63	1 (0%) 90 87	18, 29, 42, 72	0
2	C	214/237 (90%)	0.35	19 (8%) 15 13	26, 44, 102, 130	0
2	H	232/237 (97%)	-0.30	3 (1%) 75 71	24, 37, 58, 89	0
3	D	196/216 (90%)	1.14	53 (27%) 1 1	30, 68, 119, 145	0
3	L	211/216 (97%)	0.47	25 (11%) 9 7	22, 49, 93, 114	0
All	All	1627/1844 (88%)	-0.12	101 (6%) 26 23	17, 33, 92, 145	0

All (101) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
3	L	145	PRO	6.0
3	L	144	TYR	5.9
3	D	212	PRO	5.9
3	D	210	VAL	5.9
2	C	208	LEU	5.3
3	L	146	GLY	4.7
3	D	2	SER	4.6
2	C	232	PRO	4.5
3	L	32	TYR	4.5
3	D	184	LEU	4.4
3	D	32	TYR	4.4
3	D	26	SER	4.4
3	D	189	TRP	4.3
3	D	27	SER	4.0
2	C	203	VAL	4.0
3	D	30	GLY	3.8
3	D	211	ALA	3.8
3	L	2	SER	3.8
3	D	126	SER	3.8

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Mol	Chain	Res	Type	RSRZ
3	D	33	ASN	3.7
2	C	202	THR	3.7
3	L	33	ASN	3.7
3	L	27	SER	3.6
2	C	214	ILE	3.6
3	D	29	VAL	3.5
3	D	147	ALA	3.5
3	D	162	GLY	3.4
2	C	229	LYS	3.4
3	L	207	GLU	3.4
3	L	200	THR	3.4
3	L	161	ALA	3.3
3	D	127	GLU	3.2
3	D	202	GLU	3.2
2	H	110	ILE	3.1
2	C	159	CYS	3.1
3	D	28	ASP	3.1
3	L	160	LYS	3.0
3	D	145	PRO	3.0
3	D	118	SER	3.0
2	C	230	VAL	2.9
3	D	3	VAL	2.9
2	C	201	VAL	2.9
3	D	206	VAL	2.9
2	C	179	THR	2.9
3	D	196	SER	2.8
3	L	158	PRO	2.8
3	D	152	TRP	2.8
3	D	144	TYR	2.8
3	L	25	ALA	2.7
3	D	115	ALA	2.7
3	D	163	VAL	2.7
3	D	209	THR	2.7
3	D	172	SER	2.6
3	L	31	ASP	2.6
2	C	206	SER	2.6
3	L	52	ASP	2.6
3	D	52	ASP	2.6
2	H	146	SER	2.6
3	D	121	LEU	2.6
3	D	136	LEU	2.6
3	D	120	THR	2.6

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Mol	Chain	Res	Type	RSRZ
1	B	469	LYS	2.6
3	L	193	ARG	2.5
3	D	135	THR	2.5
3	L	153	LYS	2.5
3	L	29	VAL	2.5
3	D	205	THR	2.5
3	D	167	THR	2.4
3	L	115	ALA	2.4
3	L	116	ALA	2.4
3	D	116	ALA	2.4
3	D	169	SER	2.4
3	D	170	LYS	2.4
2	C	143	LEU	2.3
2	C	144	ALA	2.3
2	C	145	PRO	2.3
3	D	124	PRO	2.3
2	C	207	SER	2.3
3	L	28	ASP	2.3
3	D	117	PRO	2.3
3	D	139	LEU	2.3
3	D	150	VAL	2.2
3	D	182	LEU	2.2
3	L	26	SER	2.2
3	D	180	SER	2.2
3	D	204	SER	2.2
2	C	186	PRO	2.2
2	C	177	ALA	2.2
3	D	151	ALA	2.2
2	C	28	PRO	2.2
3	D	186	PRO	2.2
3	D	208	LYS	2.2
3	D	95	ARG	2.1
2	C	200	VAL	2.1
3	D	129	LEU	2.1
3	D	119	VAL	2.0
3	L	30	GLY	2.0
2	H	111	GLY	2.0
3	L	155	ASP	2.0
3	L	212	PRO	2.0
3	D	31	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 [i](#)

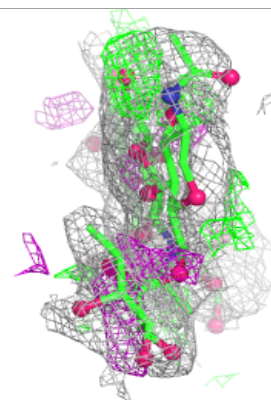
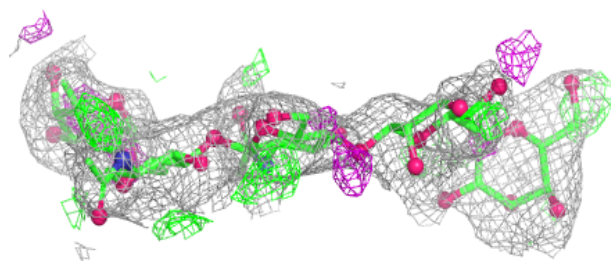
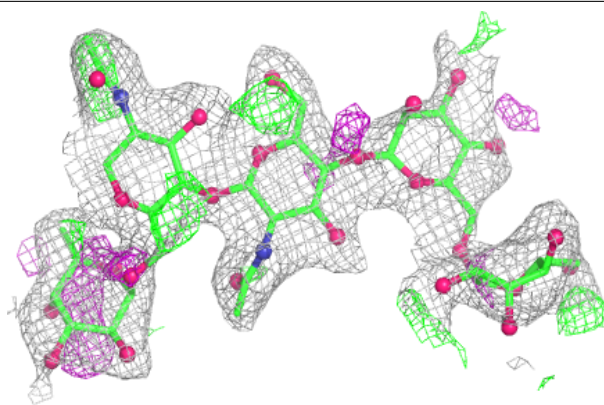
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists 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
4	NAG	E	1	14/15	-	-	52,64,69,76	0
4	NAG	E	2	14/15	-	-	46,57,70,79	0
4	BMA	E	3	11/12	-	-	96,106,112,113	0
4	MAN	E	4	11/12	-	-	78,87,90,91	0
4	FUC	E	5	10/11	-	-	44,58,62,68	0
4	NAG	F	1	14/15	-	-	42,49,60,60	0
4	NAG	F	2	14/15	-	-	38,48,53,61	0
4	BMA	F	3	11/12	-	-	87,96,108,116	0
4	MAN	F	4	11/12	-	-	87,94,100,101	0
4	FUC	F	5	10/11	-	-	44,51,55,57	0

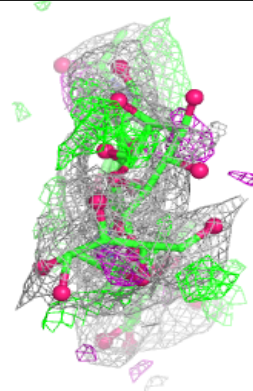
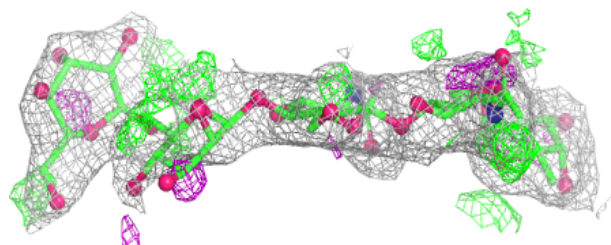
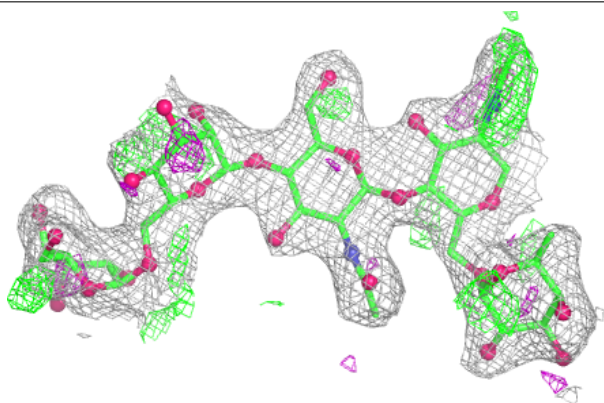
The following is a graphical depiction of the model fit to experimental electron density for oligosaccharide. Each fit is shown from different orientation to approximate a three-dimensional view.

Electron density around Chain E:

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

**Electron density around Chain F:**

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



6.4 Ligands

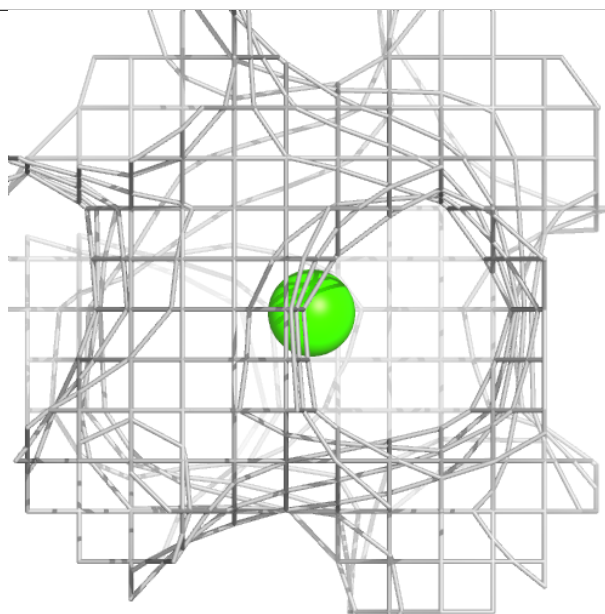
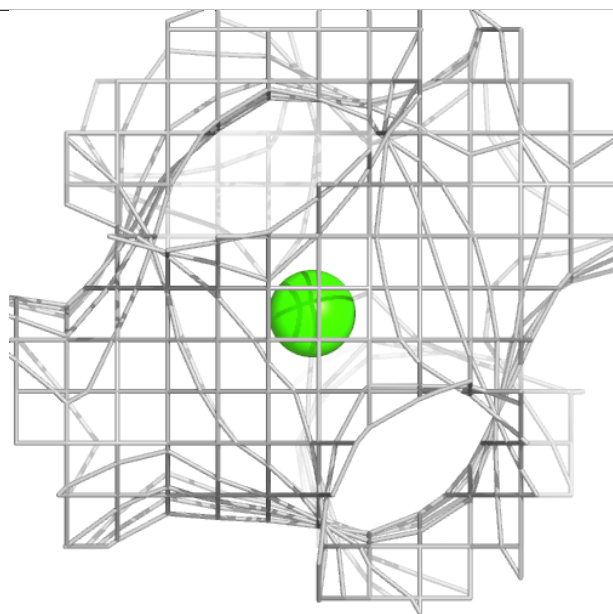
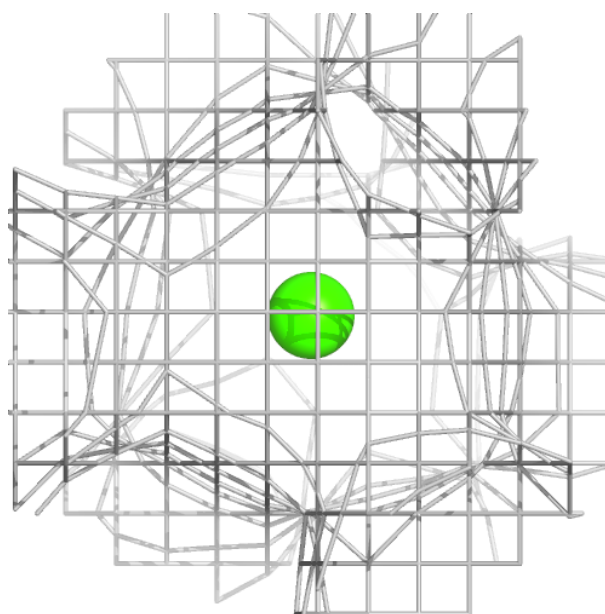
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
5	CA	A	501	1/1	0.99	0.02	27,27,27,27	0
5	CA	A	502	1/1	0.99	0.05	33,33,33,33	0
5	CA	B	501	1/1	0.99	0.03	32,32,32,32	0
5	CA	B	502	1/1	0.99	0.04	37,37,37,37	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.

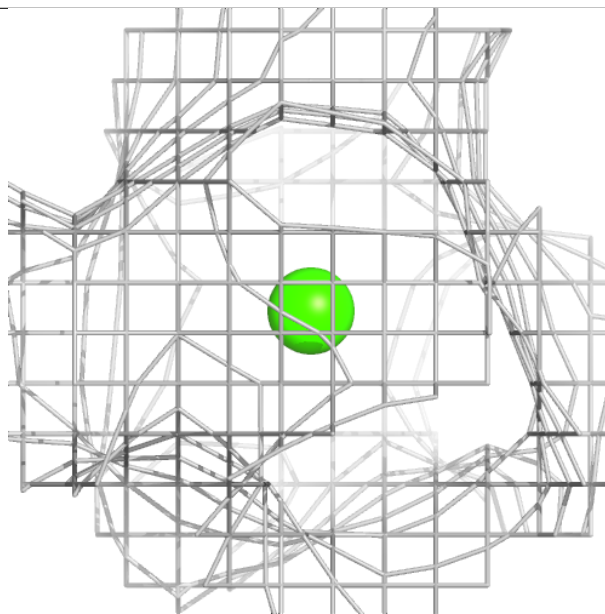
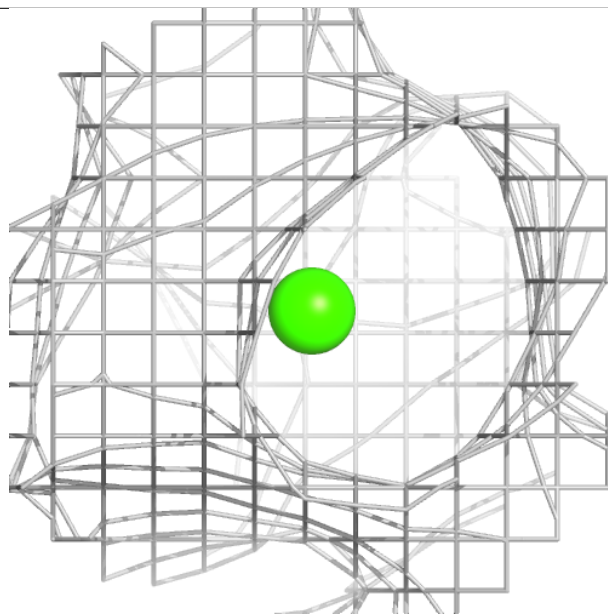
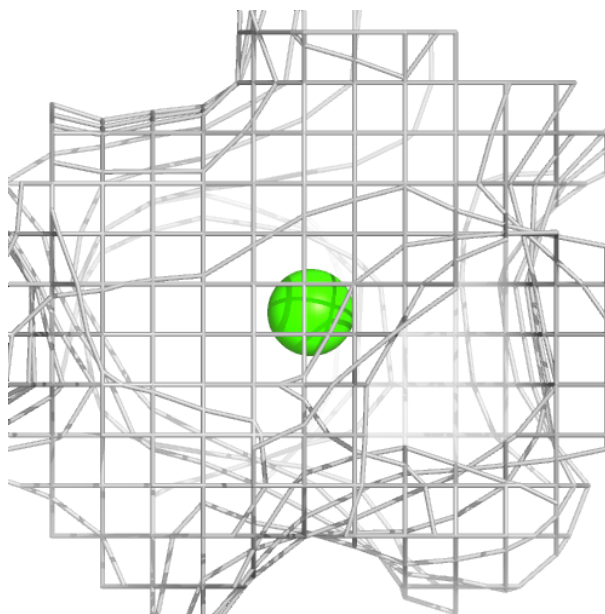
Electron density around CA A 501:

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



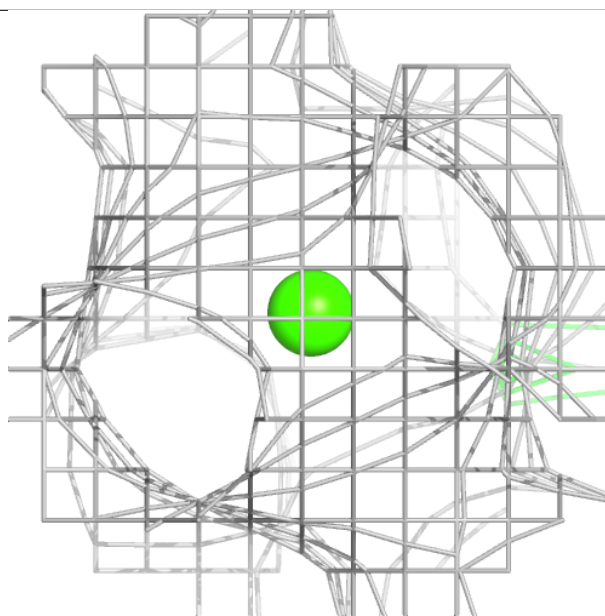
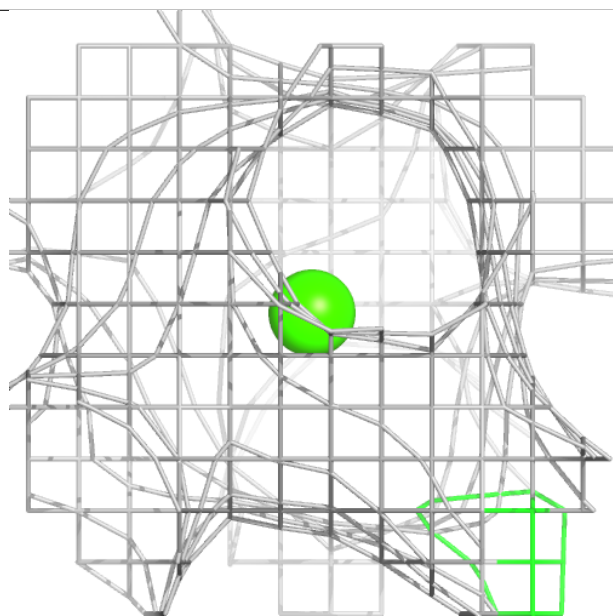
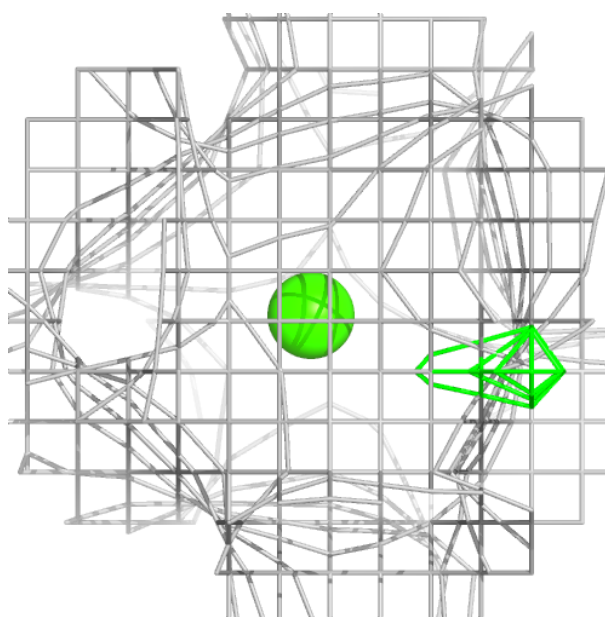
Electron density around CA A 502:

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



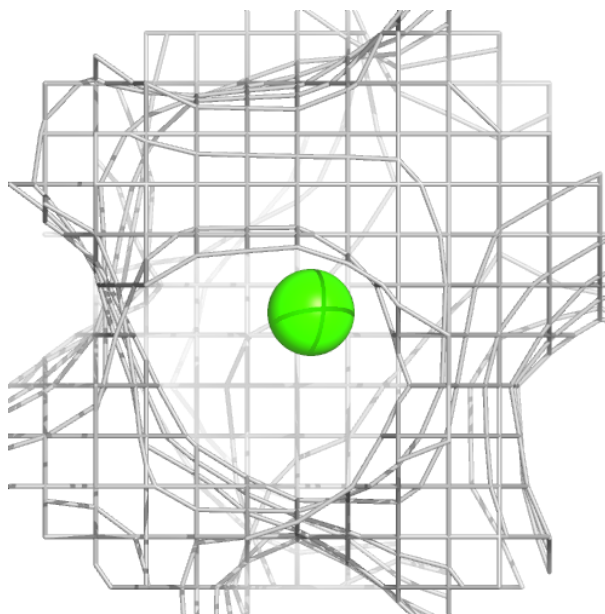
Electron density around CA B 501:

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



Electron density around CA B 502:

$2mF_o-DF_c$ (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.