



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 10:37 AM UTC

PDB ID : 2W68 / pdb_00002w68
Title : ENHANCING THE RECEPTOR AFFINITY OF THE SIALIC ACID-BINDING DOMAIN OF VIBRIO CHOLERAE SIALIDASE THROUGH MULTIVALENCY
Authors : Connaris, H.; Crocker, P.R.; Taylor, G.L.
Deposited on : 2008-12-17
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

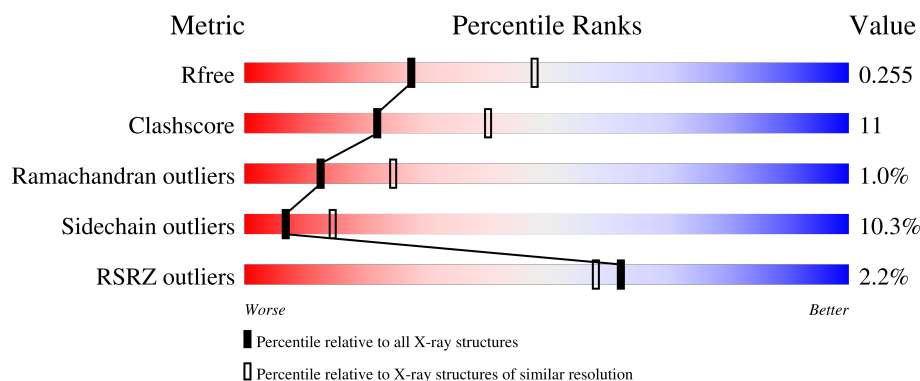
1 Overall quality at a glance

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	195	78% 17% . .
1	B	195	2% 81% 13% 5% .
1	C	195	4% 66% 28% 6% ..
2	D	3	33% 33% 33%
2	E	3	67% 33%

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Mol	Chain	Length	Quality of chain
2	F	3	 33% 67%

2 Entry composition [i](#)

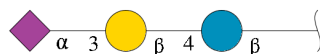
There are 4 unique types of molecules in this entry. The entry contains 4696 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 SIALIDASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	195	Total	C	N	O	S	0	0	0
			1483	927	253	296	7			
1	B	193	Total	C	N	O	S	0	0	0
			1474	922	251	294	7			
1	C	194	Total	C	N	O	S	0	0	0
			1476	924	252	293	7			

- Molecule 2 is an oligosaccharide called N-acetyl-alpha-neuraminic acid-(2-3)-beta-D-galactopyranose-(1-4)-beta-D-glucopyranose.



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	D	3	Total	C	N	O	0	0	0
			43	23	1	19			
2	E	3	Total	C	N	O	0	0	0
			43	23	1	19			
2	F	3	Total	C	N	O	0	0	0
			43	23	1	19			

- Molecule 3 is CALCIUM ION (CCD ID: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	Ca	0	0
			1	1		

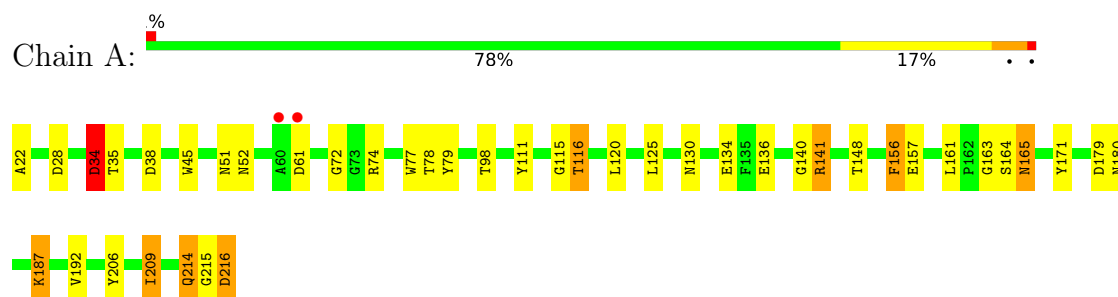
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	40	Total 40	O 40	0	0
4	B	44	Total 44	O 44	0	0
4	C	49	Total 49	O 49	0	0

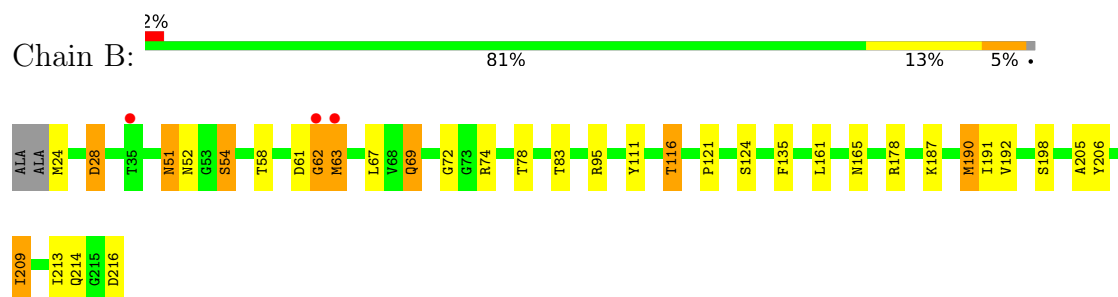
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

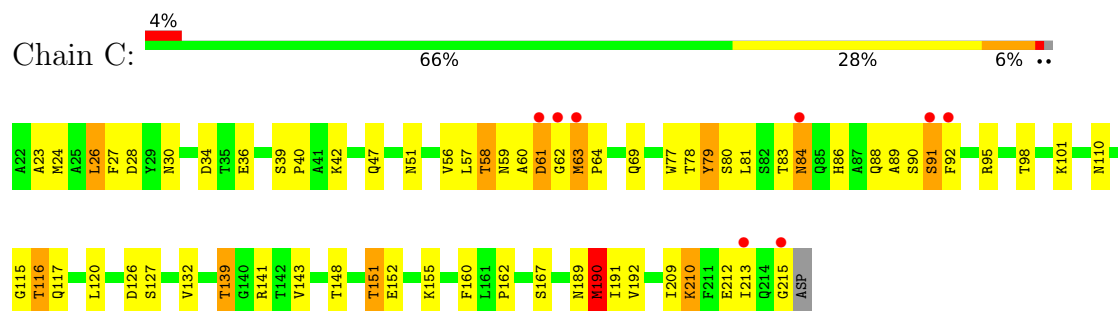
• Molecule 1: SIALIDASE



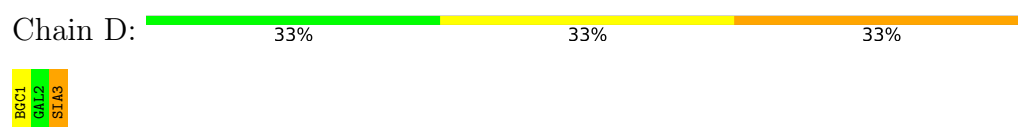
• Molecule 1: SIALIDASE



• Molecule 1: SIALIDASE



• Molecule 2: N-acetyl-alpha-neuraminic acid-(2-3)-beta-D-galactopyranose-(1-4)-beta-D-glucopyranose



- Molecule 2: N-acetyl-alpha-neuraminic acid-(2-3)-beta-D-galactopyranose-(1-4)-beta-D-glucopyranose

Chain E:  67% 33%

BGC1
GAL2
STA3

- Molecule 2: N-acetyl-alpha-neuraminic acid-(2-3)-beta-D-galactopyranose-(1-4)-beta-D-glucopyranose

Chain F:  33% 67%

BGC1
GAL2
STA3

4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	138.60Å 197.62Å 82.98Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	28.88 – 2.50 28.88 – 2.50	Depositor EDS
% Data completeness (in resolution range)	95.4 (28.88-2.50) 95.3 (28.88-2.50)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.23 (at 2.51Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.215 , 0.256 0.213 , 0.255	Depositor DCC
R_{free} test set	1911 reflections (5.06%)	wwPDB-VP
Wilson B-factor (Å ²)	47.1	Xtriage
Anisotropy	0.307	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 25.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	4696	wwPDB-VP
Average B, all atoms (Å ²)	46.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.64% 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: SIA, CA, BGC, GAL

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

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# $ Z > 5$	RMSZ	# $ Z > 5$
1	A	1.00	0/1516	1.16	11/2058 (0.5%)
1	B	0.95	0/1507	1.04	3/2046 (0.1%)
1	C	1.01	0/1509	1.15	8/2049 (0.4%)
All	All	0.99	0/4532	1.12	22/6153 (0.4%)

There are no bond length outliers.

All (22) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	51	ASN	CB-CA-C	-7.89	107.48	116.63
1	A	120	LEU	CA-C-N	-7.38	110.29	120.25
1	A	120	LEU	C-N-CA	-7.38	110.29	120.25
1	A	163	GLY	N-CA-C	-7.27	101.08	110.38
1	C	115	GLY	N-CA-C	-6.96	105.87	114.92
1	C	28	ASP	N-CA-C	6.87	118.22	108.74
1	C	63	MET	N-CA-C	6.72	114.05	108.07
1	B	111	TYR	N-CA-C	6.17	117.52	108.14
1	B	124	SER	N-CA-C	5.87	116.84	108.74
1	C	162	PRO	N-CA-C	5.80	119.78	111.13
1	B	28	ASP	N-CA-C	5.69	117.84	109.24
1	A	115	GLY	N-CA-C	-5.65	107.97	114.69
1	C	51	ASN	CB-CA-C	-5.63	110.10	116.63
1	A	214	GLN	N-CA-C	5.55	117.23	110.41
1	A	164	SER	N-CA-C	-5.53	106.20	113.17
1	A	34	ASP	CB-CA-C	-5.50	102.07	109.71
1	C	59	ASN	N-CA-C	-5.40	100.59	109.40
1	A	164	SER	CA-C-N	5.25	127.63	122.28
1	A	164	SER	C-N-CA	5.25	127.63	122.28
1	A	111	TYR	N-CA-C	5.16	116.14	108.60
1	C	120	LEU	CA-C-N	-5.09	114.75	120.14

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	120	LEU	C-N-CA	-5.09	114.75	120.14

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	1483	0	1398	24	1
1	B	1474	0	1392	28	1
1	C	1476	0	1398	44	0
2	D	43	0	37	1	0
2	E	43	0	37	2	0
2	F	43	0	37	1	0
3	A	1	0	0	0	0
4	A	40	0	0	6	0
4	B	44	0	0	10	0
4	C	49	0	0	14	1
All	All	4696	0	4299	95	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 11.

All (95) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:24:MET:HG3	4:B:2042:HOH:O	1.05	1.22
1:C:98:THR:HG22	1:C:209:ILE:HD13	1.17	1.15
1:B:74:ARG:NH1	2:E:3:SIA:O1A	1.98	0.96
1:C:98:THR:HG22	1:C:209:ILE:CD1	1.97	0.93
1:C:63:MET:HB2	1:C:64:PRO:HD2	1.55	0.86
1:B:24:MET:CG	4:B:2042:HOH:O	1.79	0.79
1:C:60:ALA:O	1:C:62:GLY:N	2.16	0.77
1:C:101:LYS:HE3	1:C:151:THR:O	1.87	0.74
1:C:80:SER:O	1:C:81:LEU:HD23	1.88	0.73

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:190:MET:HG3	1:B:192:VAL:CG2	2.20	0.70
1:A:157:GLU:HB2	1:A:171:TYR:HB2	1.75	0.67
1:C:95:ARG:NH1	1:C:212:GLU:OE1	2.30	0.64
1:C:57:LEU:HD13	1:C:58:THR:N	2.13	0.64
1:B:190:MET:HG3	1:B:192:VAL:HG23	1.79	0.64
1:C:63:MET:HB2	1:C:64:PRO:CD	2.28	0.63
1:B:24:MET:N	4:B:2001:HOH:O	2.30	0.63
1:A:165:ASN:HB2	4:A:2033:HOH:O	1.98	0.62
1:C:26:LEU:HD12	1:C:27:PHE:N	2.14	0.62
1:B:54:SER:HB3	1:B:69:GLN:OE1	1.99	0.62
1:A:187:LYS:N	1:A:187:LYS:HD3	2.15	0.61
1:A:136:GLU:OE2	1:C:139:THR:HG23	2.00	0.61
1:B:63:MET:SD	1:B:63:MET:N	2.73	0.61
1:C:23:ALA:O	1:C:215:GLY:C	2.44	0.60
1:B:52:ASN:HB2	1:B:72:GLY:O	2.01	0.60
1:C:42:LYS:HE2	4:C:2005:HOH:O	2.01	0.59
1:C:160:PHE:CG	4:C:2037:HOH:O	2.52	0.58
1:A:34:ASP:O	1:A:38:ASP:HB3	2.04	0.57
1:B:74:ARG:HD3	2:E:3:SIA:O1A	2.04	0.57
2:F:1:BGC:H4	2:F:2:GAL:H61	1.85	0.57
1:B:206:TYR:CD2	1:B:209:ILE:HD11	2.40	0.57
1:C:79:TYR:HB3	1:C:191:ILE:HD12	1.86	0.56
1:C:95:ARG:HD3	4:C:2021:HOH:O	2.05	0.56
1:C:116:THR:HG22	1:C:117:GLN:HG2	1.86	0.56
1:A:98:THR:HG23	1:A:209:ILE:HD13	1.86	0.56
1:C:78:THR:HG22	1:C:192:VAL:HG22	1.89	0.55
1:B:24:MET:CB	4:B:2042:HOH:O	2.37	0.54
1:C:60:ALA:C	1:C:62:GLY:N	2.67	0.53
1:A:206:TYR:CD2	1:A:209:ILE:HD11	2.44	0.53
1:C:36:GLU:HG3	4:C:2004:HOH:O	2.10	0.52
1:B:216:ASP:HA	4:B:2043:HOH:O	2.09	0.52
1:A:136:GLU:OE2	1:C:139:THR:CG2	2.58	0.51
1:B:121:PRO:HB3	1:B:135:PHE:CE2	2.46	0.51
1:B:213:ILE:HG22	1:B:214:GLN:O	2.11	0.51
1:B:95:ARG:HD2	4:B:2041:HOH:O	2.09	0.51
1:B:165:ASN:ND2	4:B:2032:HOH:O	2.44	0.50
1:C:84:ASN:N	1:C:84:ASN:OD1	2.44	0.50
1:C:80:SER:C	1:C:81:LEU:HD23	2.37	0.50
1:B:62:GLY:C	1:B:63:MET:SD	2.96	0.49
1:A:116:THR:CG2	4:A:2023:HOH:O	2.61	0.49
1:B:24:MET:HB2	4:B:2042:HOH:O	2.08	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:213:ILE:O	1:C:213:ILE:HG22	2.13	0.48
1:C:189:ASN:O	1:C:190:MET:HB3	2.14	0.48
1:C:132:VAL:HG12	1:C:143:VAL:HA	1.96	0.48
1:B:61:ASP:C	1:B:63:MET:H	2.22	0.47
1:B:28:ASP:HB2	1:B:209:ILE:O	2.15	0.47
1:A:116:THR:HG22	4:A:2023:HOH:O	2.13	0.47
1:C:189:ASN:HB3	4:C:2045:HOH:O	2.14	0.47
1:C:57:LEU:HD13	1:C:58:THR:H	1.79	0.47
1:A:216:ASP:N	1:A:216:ASP:OD1	2.49	0.46
1:C:40:PRO:HG2	1:C:77:TRP:CE2	2.50	0.46
1:C:86:HIS:O	1:C:89:ALA:HB3	2.16	0.46
1:C:152:GLU:HG3	4:C:2033:HOH:O	2.16	0.46
1:A:125:LEU:HA	1:A:130:ASN:O	2.15	0.46
1:B:95:ARG:CD	4:B:2041:HOH:O	2.64	0.46
1:C:160:PHE:CD2	4:C:2037:HOH:O	2.67	0.46
1:A:179:ASP:O	1:A:180:ASN:HB2	2.16	0.46
1:C:62:GLY:HA2	4:C:2009:HOH:O	2.15	0.46
1:C:84:ASN:ND2	4:C:2016:HOH:O	2.49	0.45
1:A:215:GLY:HA2	4:A:2001:HOH:O	2.16	0.45
1:A:52:ASN:HB2	1:A:72:GLY:O	2.16	0.45
1:A:78:THR:HG22	1:A:192:VAL:HG22	1.97	0.45
1:C:98:THR:CG2	1:C:209:ILE:CD1	2.83	0.45
1:B:78:THR:HA	1:B:191:ILE:O	2.17	0.45
1:A:74:ARG:NH1	2:D:3:SIA:O1B	2.47	0.44
1:C:26:LEU:HD11	1:C:210:LYS:HG3	2.00	0.43
1:A:134:GLU:HG2	1:A:140:GLY:O	2.18	0.43
1:C:24:MET:HE1	4:C:2047:HOH:O	2.18	0.43
1:B:24:MET:HE3	4:B:2043:HOH:O	2.19	0.43
1:B:67:LEU:HD12	1:B:205:ALA:HB2	2.01	0.43
1:C:126:ASP:OD1	1:C:126:ASP:C	2.63	0.42
1:C:110:ASN:HA	1:C:192:VAL:O	2.19	0.42
1:C:190:MET:HG3	1:C:192:VAL:HG23	2.02	0.42
1:B:69:GLN:O	1:B:69:GLN:NE2	2.52	0.42
1:B:178:ARG:HE	1:B:178:ARG:HB2	1.27	0.41
1:C:62:GLY:CA	4:C:2009:HOH:O	2.68	0.41
1:A:165:ASN:CB	4:A:2033:HOH:O	2.65	0.41
1:B:51:ASN:HB2	1:B:52:ASN:H	1.67	0.41
1:A:45:TRP:CE3	1:A:79:TYR:HB2	2.56	0.41
1:A:77:TRP:NE1	4:A:2010:HOH:O	2.54	0.41
1:A:156:PHE:N	1:A:156:PHE:CD1	2.88	0.41
1:A:45:TRP:CZ3	1:A:79:TYR:HB2	2.56	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:141:ARG:HD3	4:C:2048:HOH:O	2.21	0.40
1:C:88:GLN:HA	1:C:91:SER:HB3	2.04	0.40
1:C:190:MET:CE	4:C:2013:HOH:O	2.69	0.40
1:C:24:MET:HG3	4:C:2001:HOH:O	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)
1:B:24:MET:CE	4:C:2002:HOH:O[3_555]	1.98	0.22
1:A:22:ALA:N	1:A:28:ASP:N[3_554]	2.01	0.19

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	193/195 (99%)	181 (94%)	12 (6%)	0	100	100
1	B	191/195 (98%)	178 (93%)	11 (6%)	2 (1%)	12	24
1	C	192/195 (98%)	181 (94%)	7 (4%)	4 (2%)	5	9
All	All	576/585 (98%)	540 (94%)	30 (5%)	6 (1%)	12	24

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	61	ASP
1	B	62	GLY
1	B	116	THR
1	C	91	SER
1	C	34	ASP
1	C	190	MET

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	155/155 (100%)	142 (92%)	13 (8%)	10	22
1	B	155/155 (100%)	143 (92%)	12 (8%)	12	25
1	C	154/155 (99%)	131 (85%)	23 (15%)	3	6
All	All	464/465 (100%)	416 (90%)	48 (10%)	7	14

All (48) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	34	ASP
1	A	35	THR
1	A	61	ASP
1	A	116	THR
1	A	141	ARG
1	A	148	THR
1	A	156	PHE
1	A	161	LEU
1	A	165	ASN
1	A	187	LYS
1	A	209	ILE
1	A	214	GLN
1	A	216	ASP
1	B	51	ASN
1	B	54	SER
1	B	58	THR
1	B	63	MET
1	B	69	GLN
1	B	83	THR
1	B	116	THR
1	B	161	LEU
1	B	187	LYS
1	B	190	MET
1	B	198	SER
1	B	209	ILE
1	C	26	LEU

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Mol	Chain	Res	Type
1	C	30	ASN
1	C	39	SER
1	C	47	GLN
1	C	56	VAL
1	C	58	THR
1	C	61	ASP
1	C	69	GLN
1	C	79	TYR
1	C	83	THR
1	C	84	ASN
1	C	90	SER
1	C	92	PHE
1	C	116	THR
1	C	127	SER
1	C	139	THR
1	C	141	ARG
1	C	148	THR
1	C	151	THR
1	C	155	LYS
1	C	167	SER
1	C	190	MET
1	C	210	LYS

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

Mol	Chain	Res	Type
1	A	69	GLN
1	A	165	ASN
1	B	69	GLN
1	B	84	ASN
1	B	165	ASN
1	B	182	GLN
1	B	195	ASN
1	C	30	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

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
2	BGC	D	1	2	12,12,12	0.66	0	17,17,17	1.16	3 (17%)
2	GAL	D	2	2	11,11,12	0.33	0	15,15,17	0.89	0
2	SIA	D	3	2	20,20,21	0.79	0	21,28,31	1.45	3 (14%)
2	BGC	E	1	2	12,12,12	0.65	0	17,17,17	1.70	3 (17%)
2	GAL	E	2	2	11,11,12	0.46	0	15,15,17	1.51	4 (26%)
2	SIA	E	3	2	20,20,21	1.56	1 (5%)	21,28,31	1.75	5 (23%)
2	BGC	F	1	2	12,12,12	0.79	0	17,17,17	1.17	2 (11%)
2	GAL	F	2	2	11,11,12	0.62	0	15,15,17	2.50	5 (33%)
2	SIA	F	3	2	20,20,21	1.18	1 (5%)	21,28,31	1.66	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
2	BGC	D	1	2	-	2/2/22/22	0/1/1/1
2	GAL	D	2	2	-	2/2/19/22	0/1/1/1
2	SIA	D	3	2	-	2/18/34/38	0/1/1/1
2	BGC	E	1	2	-	2/2/22/22	0/1/1/1
2	GAL	E	2	2	-	2/2/19/22	0/1/1/1
2	SIA	E	3	2	-	2/18/34/38	0/1/1/1
2	BGC	F	1	2	-	2/2/22/22	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	GAL	F	2	2	-	2/2/19/22	0/1/1/1
2	SIA	F	3	2	-	2/18/34/38	0/1/1/1

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	E	3	SIA	C2-C1	5.89	1.59	1.52
2	F	3	SIA	C2-C1	3.52	1.56	1.52

All (28) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	F	2	GAL	C1-O5-C5	6.17	120.45	112.19
2	F	2	GAL	C1-C2-C3	-5.07	102.26	109.64
2	E	3	SIA	C6-C5-N5	-4.39	103.90	110.91
2	F	3	SIA	O6-C2-C3	-4.18	104.94	110.56
2	F	3	SIA	C8-C7-C6	-4.09	105.37	113.05
2	E	1	BGC	O5-C5-C4	3.81	116.56	109.70
2	D	3	SIA	C4-C5-N5	-3.48	103.58	110.44
2	E	3	SIA	C8-C7-C6	-3.45	106.56	113.05
2	E	1	BGC	C1-O5-C5	3.41	120.25	113.65
2	E	1	BGC	C3-C4-C5	3.19	116.01	110.23
2	E	2	GAL	C1-C2-C3	-3.17	105.03	109.64
2	F	1	BGC	O5-C5-C6	3.11	114.16	106.44
2	F	2	GAL	C2-C3-C4	-2.74	106.05	110.86
2	F	3	SIA	C6-C5-N5	-2.68	106.63	110.91
2	E	3	SIA	O6-C2-C1	2.63	112.68	107.72
2	E	3	SIA	C4-C5-N5	-2.43	105.65	110.44
2	D	3	SIA	O1B-C1-C2	2.40	118.94	112.71
2	E	2	GAL	O3-C3-C2	-2.35	105.27	110.05
2	F	1	BGC	C3-C4-C5	2.30	114.40	110.23
2	F	2	GAL	O3-C3-C4	2.29	115.78	110.38
2	E	2	GAL	O5-C5-C6	2.20	111.94	107.66
2	E	3	SIA	C11-C10-N5	-2.18	112.50	116.12
2	D	3	SIA	C6-C5-N5	-2.14	107.49	110.91
2	E	2	GAL	O5-C5-C4	-2.14	105.63	110.83
2	D	1	BGC	O5-C1-C2	2.12	114.02	110.30
2	F	2	GAL	O6-C6-C5	-2.04	104.38	111.33
2	D	1	BGC	C1-O5-C5	2.01	117.55	113.65
2	D	1	BGC	C3-C4-C5	2.01	113.87	110.23

There are no chirality outliers.

All (18) torsion outliers are listed below:

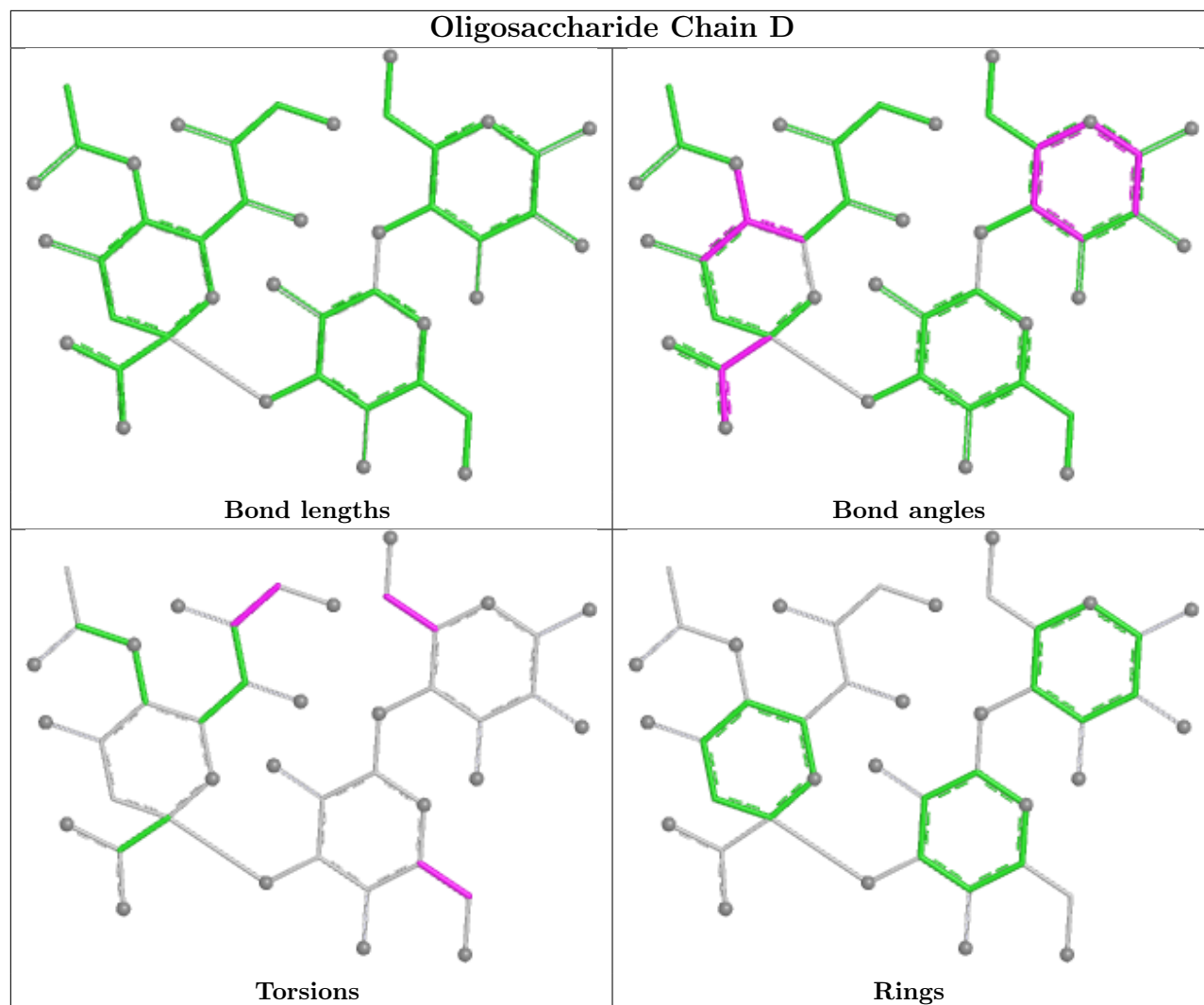
Mol	Chain	Res	Type	Atoms
2	D	3	SIA	O8-C8-C9-O9
2	E	3	SIA	C7-C8-C9-O9
2	E	3	SIA	O8-C8-C9-O9
2	F	3	SIA	O8-C8-C9-O9
2	E	2	GAL	O5-C5-C6-O6
2	D	3	SIA	C7-C8-C9-O9
2	F	3	SIA	C7-C8-C9-O9
2	E	2	GAL	C4-C5-C6-O6
2	E	1	BGC	C4-C5-C6-O6
2	F	1	BGC	O5-C5-C6-O6
2	F	1	BGC	C4-C5-C6-O6
2	D	1	BGC	O5-C5-C6-O6
2	D	2	GAL	O5-C5-C6-O6
2	F	2	GAL	O5-C5-C6-O6
2	D	1	BGC	C4-C5-C6-O6
2	E	1	BGC	O5-C5-C6-O6
2	D	2	GAL	C4-C5-C6-O6
2	F	2	GAL	C4-C5-C6-O6

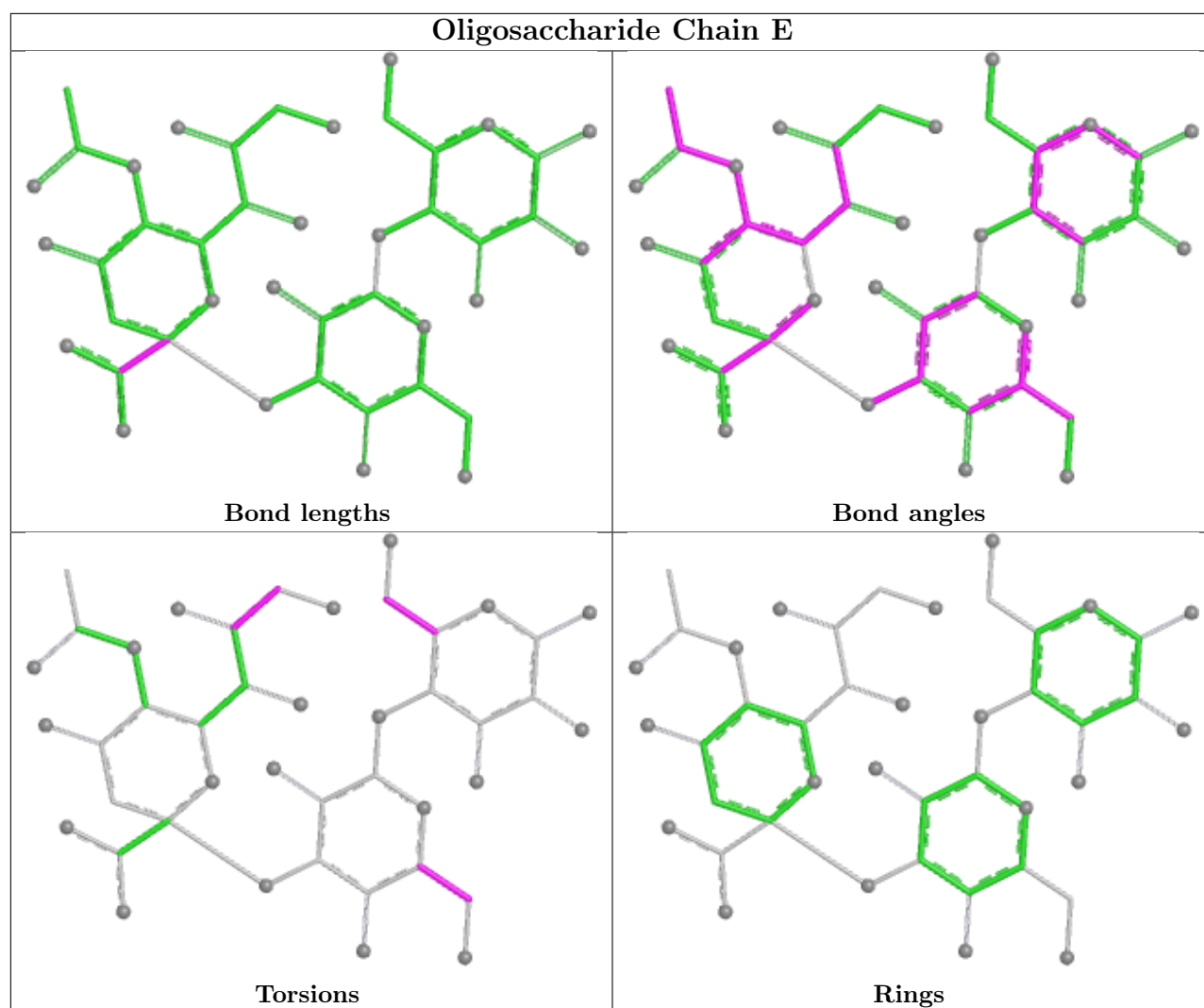
There are no ring outliers.

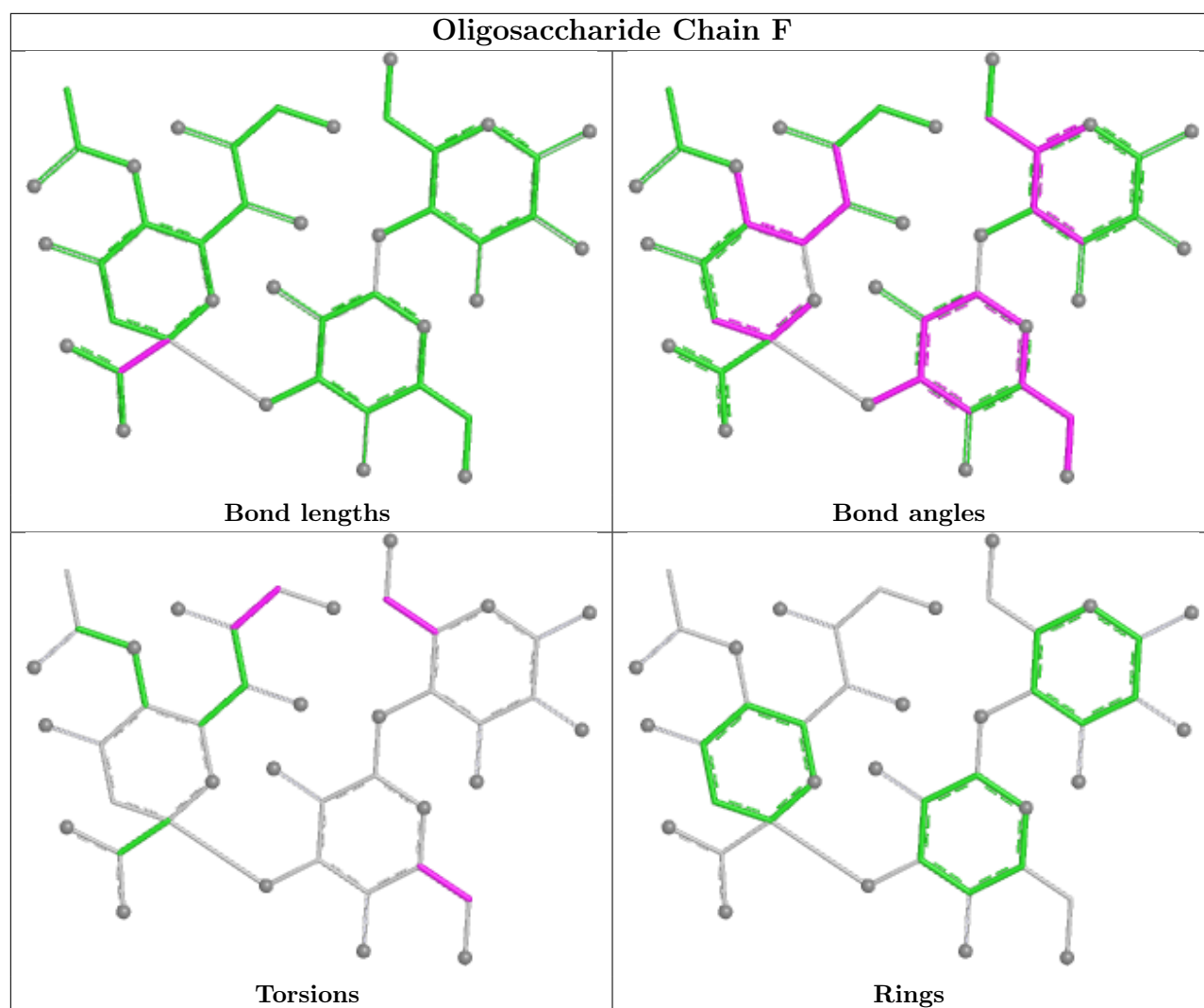
4 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	F	1	BGC	1	0
2	F	2	GAL	1	0
2	D	3	SIA	1	0
2	E	3	SIA	2	0

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







5.6 Ligand geometry [i](#)

Of 1 ligands modelled in this entry, 1 is 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 ⓘ

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	195/195 (100%)	-0.23	2 (1%) 79 76	28, 42, 57, 67	0
1	B	193/195 (98%)	-0.20	3 (1%) 70 67	29, 44, 62, 73	0
1	C	194/195 (99%)	0.30	8 (4%) 41 37	33, 51, 64, 70	0
All	All	582/585 (99%)	-0.04	13 (2%) 62 58	28, 45, 62, 73	0

All (13) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	61	ASP	3.8
1	A	60	ALA	3.7
1	C	215	GLY	3.6
1	C	92	PHE	3.5
1	C	62	GLY	3.4
1	C	213	ILE	2.9
1	A	61	ASP	2.5
1	B	35	THR	2.5
1	B	62	GLY	2.5
1	B	63	MET	2.3
1	C	91	SER	2.3
1	C	63	MET	2.2
1	C	84	ASN	2.2

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

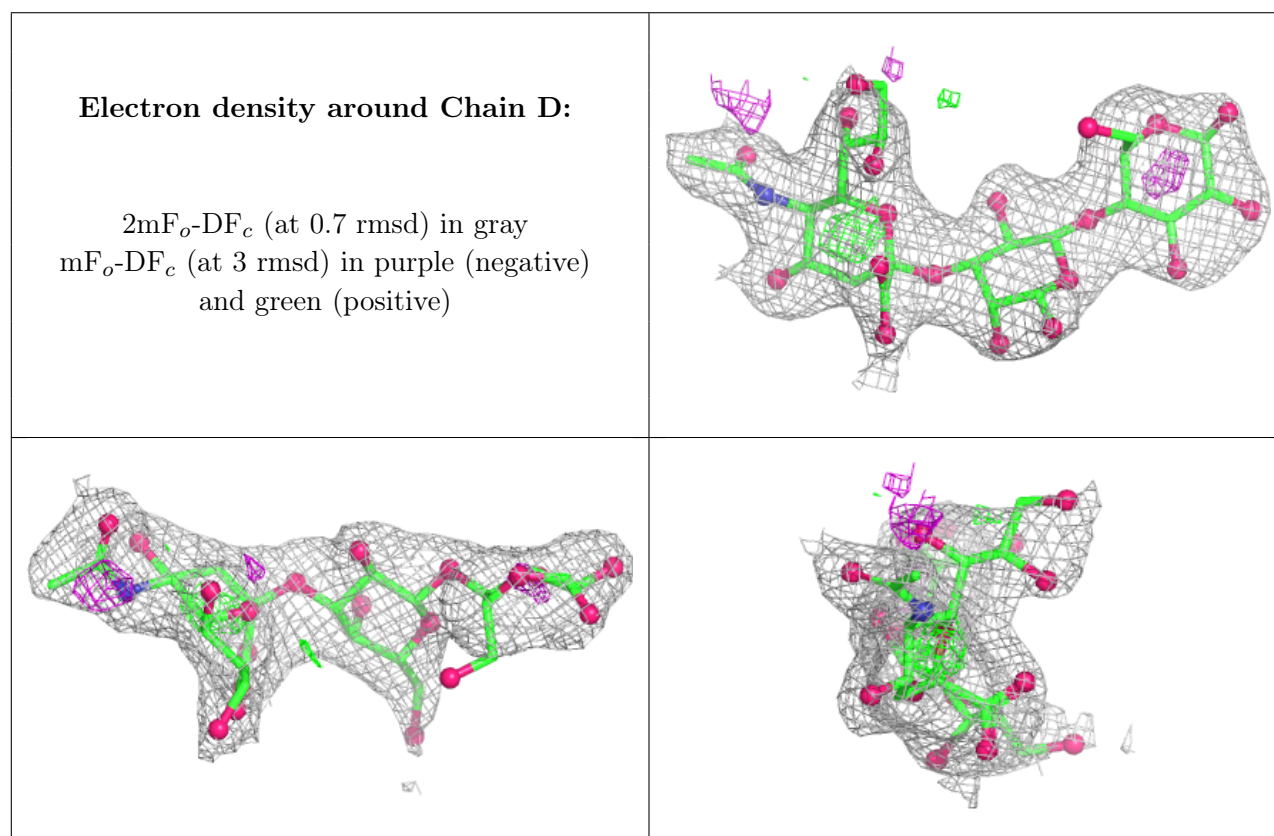
There are no non-standard protein/DNA/RNA residues in this entry.

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.

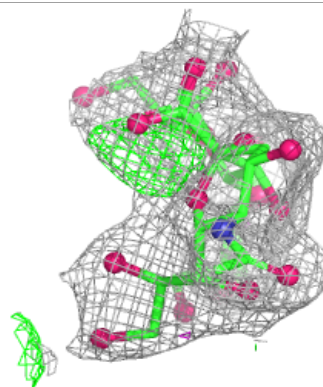
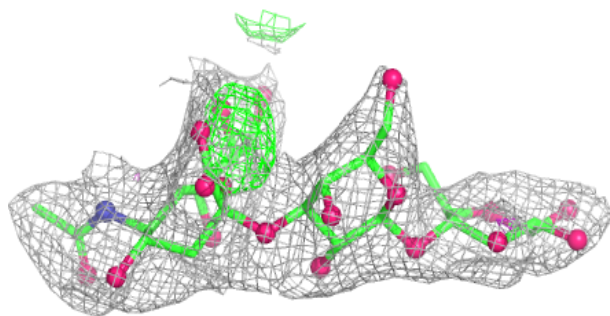
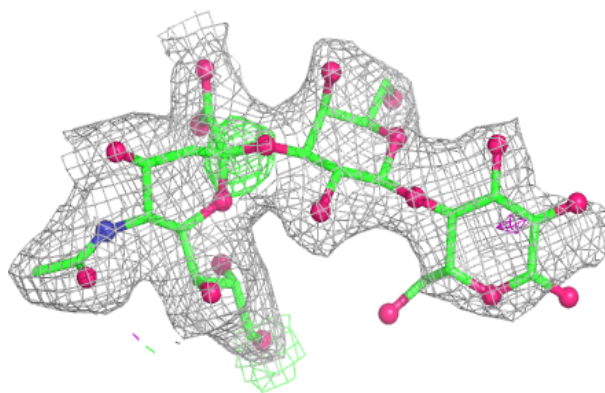
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	BGC	F	1	12/12	0.75	0.15	85,88,89,90	0
2	BGC	D	1	12/12	0.76	0.14	81,86,88,89	0
2	BGC	E	1	12/12	0.78	0.17	77,84,86,86	0
2	GAL	F	2	11/12	0.82	0.13	63,76,80,80	0
2	GAL	D	2	11/12	0.90	0.09	63,72,76,77	0
2	GAL	E	2	11/12	0.92	0.09	60,69,72,72	0
2	SIA	E	3	20/21	0.93	0.10	42,49,59,59	0
2	SIA	D	3	20/21	0.94	0.11	44,53,61,62	0
2	SIA	F	3	20/21	0.94	0.09	43,51,59,59	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.

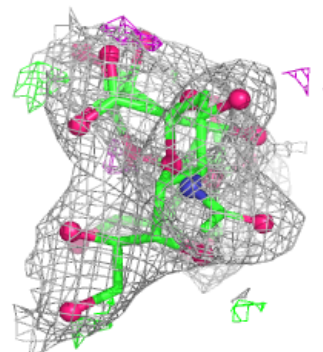
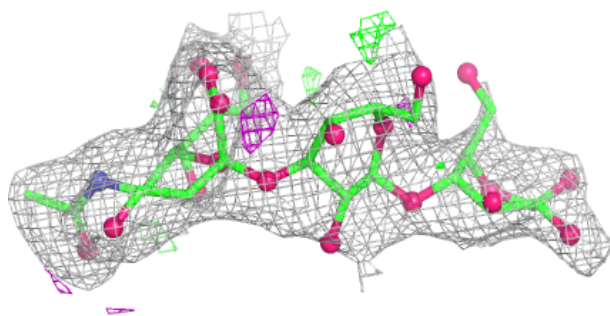
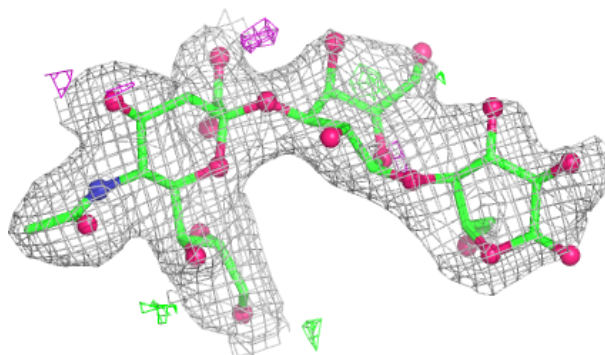


Electron density around Chain E:

$2mF_o-DF_c$ (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 [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
3	CA	A	1211	1/1	0.99	0.20	32,32,32,32	1

6.5 Other polymers [i](#)

There are no such residues in this entry.