



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 11:21 AM UTC

PDB ID : 5DNF / pdb_00005dnf
Title : Crystal structure of CC chemokine 5 (CCL5) oligomer in complex with heparin
Authors : Liang, W.G.; Tang, W.
Deposited on : 2015-09-10
Resolution : 2.55 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)
Xtrriage (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

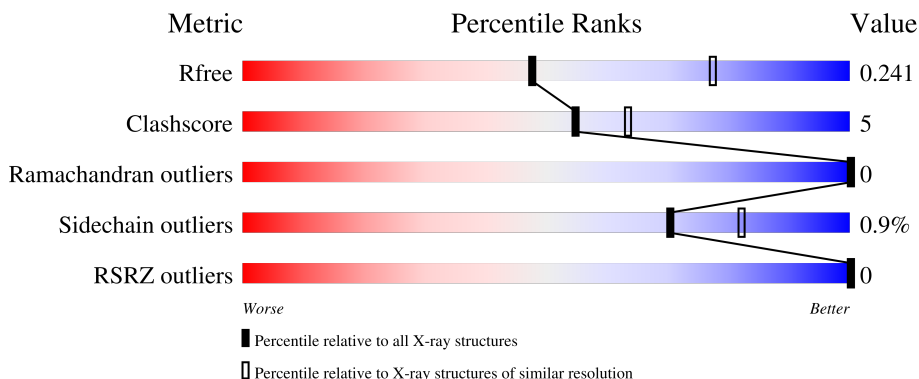
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION


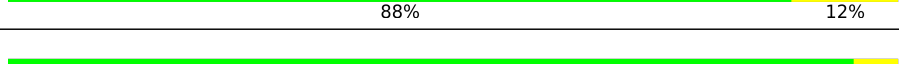
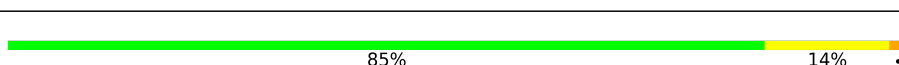

The reported resolution of this entry is 2.55 Å.

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	1091 (2.54-2.54)
Clashscore	190562	1120 (2.54-2.54)
Ramachandran outliers	187476	1106 (2.54-2.54)
Sidechain outliers	187428	1106 (2.54-2.54)
RSRZ outliers	180081	1091 (2.54-2.54)

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	65	 85% 15%
1	B	65	 92% 8%
1	C	65	 88% 12%
1	D	65	 95% 5%
1	E	65	 85% 14% .

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Mol	Chain	Length	Quality of chain
1	F	65	 88% 12%
1	G	65	 85% 15%
1	H	65	 85% 15%
1	I	65	 94% 6%
2	J	3	 100%
2	K	3	 100%
2	L	3	 67% 33%

2 Entry composition i

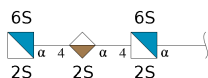
There are 8 unique types of molecules in this entry. The entry contains 5295 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 C-C motif chemokine 5.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	65	Total 525	C 333	N 93	O 94	S 5	0	0	0
1	B	65	Total 544	C 343	N 99	O 97	S 5	0	3	0
1	C	65	Total 525	C 333	N 93	O 94	S 5	0	0	0
1	D	65	Total 525	C 333	N 93	O 94	S 5	0	0	0
1	E	65	Total 525	C 333	N 93	O 94	S 5	0	1	0
1	F	65	Total 525	C 333	N 93	O 94	S 5	0	0	0
1	G	65	Total 525	C 333	N 93	O 94	S 5	0	1	0
1	H	65	Total 525	C 333	N 93	O 94	S 5	0	0	0
1	I	65	Total 525	C 333	N 93	O 94	S 5	0	0	0

- Molecule 2 is an oligosaccharide called 2-deoxy-6-O-sulfo-2-(sulfoamino)-alpha-D-glucopyranose-(1-4)-2-O-sulfo-alpha-L-idopyranuronic acid-(1-4)-2-deoxy-6-O-sulfo-2-(sulfoamino)-alpha-D-glucopyranose.



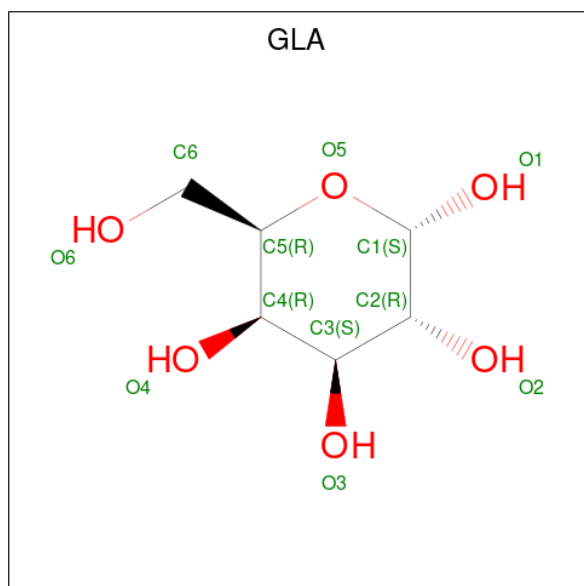
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	J	3	Total 54	C 18	N 2	O 29	S 5	0	3	0
2	K	3	Total 54	C 18	N 2	O 29	S 5	0	3	0

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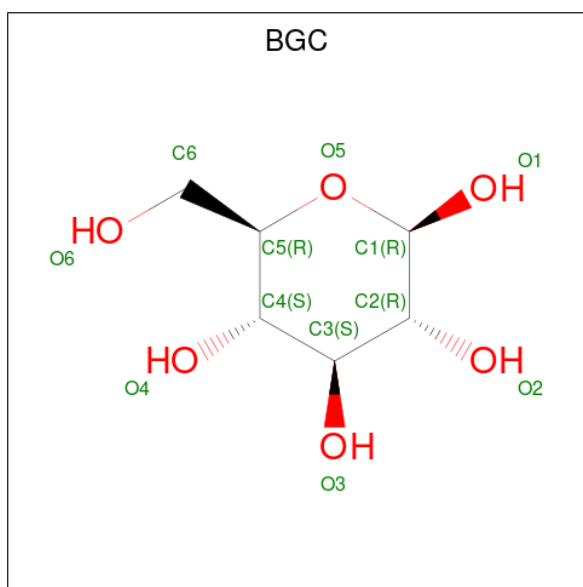
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace	
			Total	C	N	O				S
2	L	3	54	18	2	29	5	0	3	0

- Molecule 3 is alpha-D-galactopyranose (CCD ID: GLA) (formula: C₆H₁₂O₆).



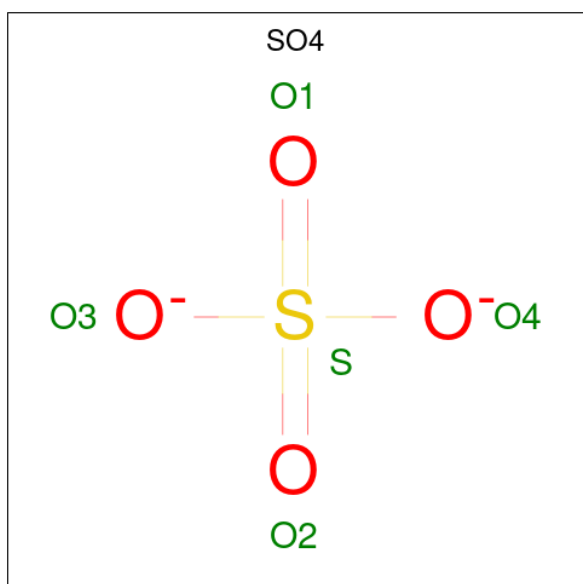
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	C	O		
3	A	1	11	6	5	0	0
3	C	1	11	6	5	0	0
3	D	1	11	6	5	0	0
3	F	1	11	6	5	0	0
3	H	1	11	6	5	0	0
3	I	1	11	6	5	0	0

- Molecule 4 is beta-D-glucopyranose (CCD ID: BGC) (formula: C₆H₁₂O₆).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	1	Total C O 12 6 6	0	0
4	C	1	Total C O 12 6 6	0	0
4	E	1	Total C O 12 6 6	0	0
4	G	1	Total C O 12 6 6	0	0
4	H	1	Total C O 12 6 6	0	0

- Molecule 5 is SULFATE ION (CCD ID: SO4) (formula: O₄S).

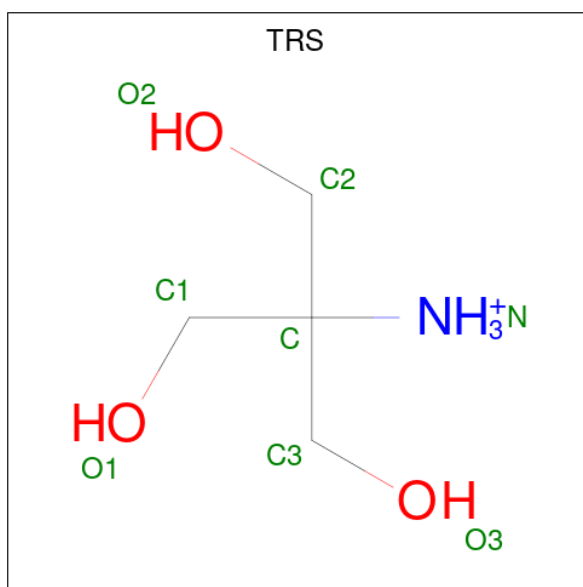


Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	1	Total O S 5 4 1	0	0
5	E	1	Total O S 5 4 1	0	0
5	F	1	Total O S 5 4 1	0	0
5	G	1	Total O S 5 4 1	0	0
5	H	1	Total O S 5 4 1	0	0

- Molecule 6 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	A	2	Total Cl 2 2	0	0
6	B	2	Total Cl 2 2	0	0
6	C	4	Total Cl 4 4	0	0
6	D	2	Total Cl 2 2	0	0
6	E	1	Total Cl 1 1	0	0
6	F	3	Total Cl 3 3	0	0
6	G	2	Total Cl 2 2	0	0
6	H	2	Total Cl 2 2	0	0
6	I	2	Total Cl 2 2	0	0

- Molecule 7 is 2-AMINO-2-HYDROXYMETHYL-PROPANE-1,3-DIOL (CCD ID: TRS) (formula: C₄H₁₂NO₃).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	
			Total	C	N			O
7	B	1	8	4	1	3	0	0


- Molecule 8 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	A	24	Total	O	0	0
			24	24		
8	B	20	Total	O	0	0
			20	20		
8	C	21	Total	O	0	0
			21	21		
8	D	19	Total	O	0	0
			19	19		
8	E	32	Total	O	0	0
			32	32		
8	F	19	Total	O	0	0
			19	19		
8	G	28	Total	O	0	0
			28	28		
8	H	24	Total	O	0	0
			24	24		
8	I	23	Total	O	0	0
			23	23		

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.

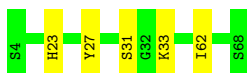
- Molecule 1: C-C motif chemokine 5

Chain A:  85% 15%




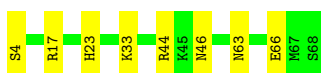
- Molecule 1: C-C motif chemokine 5

Chain B:  92% 8%



- Molecule 1: C-C motif chemokine 5

Chain C:  88% 12%




- Molecule 1: C-C motif chemokine 5

Chain D:  95% 5%




- Molecule 1: C-C motif chemokine 5

Chain E:  85% 14% .

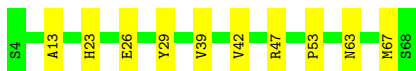
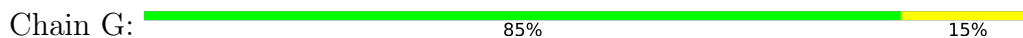


- Molecule 1: C-C motif chemokine 5

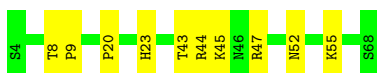
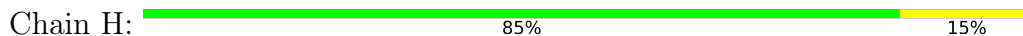
Chain F:  88% 12%



- Molecule 1: C-C motif chemokine 5



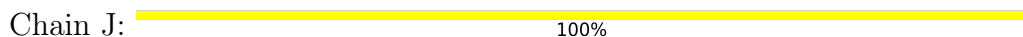
- Molecule 1: C-C motif chemokine 5



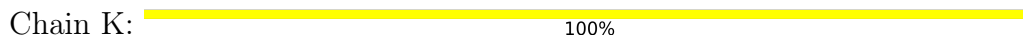
- Molecule 1: C-C motif chemokine 5



- Molecule 2: 2-deoxy-6-O-sulfo-2-(sulfoamino)-alpha-D-glucopyranose-(1-4)-2-O-sulfo-alpha-L-idopyranuronic acid-(1-4)-2-deoxy-6-O-sulfo-2-(sulfoamino)-alpha-D-glucopyranose



- Molecule 2: 2-deoxy-6-O-sulfo-2-(sulfoamino)-alpha-D-glucopyranose-(1-4)-2-O-sulfo-alpha-L-idopyranuronic acid-(1-4)-2-deoxy-6-O-sulfo-2-(sulfoamino)-alpha-D-glucopyranose



- Molecule 2: 2-deoxy-6-O-sulfo-2-(sulfoamino)-alpha-D-glucopyranose-(1-4)-2-O-sulfo-alpha-L-idopyranuronic acid-(1-4)-2-deoxy-6-O-sulfo-2-(sulfoamino)-alpha-D-glucopyranose



4 Data and refinement statistics i

Property	Value	Source
Space group	C 2 2 2	Depositor
Cell constants a, b, c, α , β , γ	121.19Å 210.06Å 123.83Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	39.70 – 2.55 39.70 – 2.55	Depositor EDS
% Data completeness (in resolution range)	96.9 (39.70-2.55) 94.3 (39.70-2.55)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	0.11	Depositor
$\langle I/\sigma(I) \rangle$ ¹	5.60 (at 2.54Å)	Xtrriage
Refinement program	PHENIX 1.9_1692	Depositor
R, R_{free}	0.180 , 0.234 0.198 , 0.241	Depositor DCC
R_{free} test set	2004 reflections (3.90%)	wwPDB-VP
Wilson B-factor (Å ²)	35.5	Xtrriage
Anisotropy	0.068	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.39 , 33.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtrriage
Estimated twinning fraction	0.488 for 1/2*h-1/2*k,-3/2*h-1/2*k,-l 0.488 for 1/2*h+1/2*k,3/2*h-1/2*k,-l	Xtrriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	5295	wwPDB-VP
Average B, all atoms (Å ²)	42.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.86% 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: SGN, GLA, IDS, SO4, CL, BGC, TRS

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	0.29	0/539	0.73	0/730
1	B	0.28	0/558	0.71	0/755
1	C	0.29	0/539	0.77	0/730
1	D	0.29	0/539	0.76	0/730
1	E	0.36	0/539	0.87	3/730 (0.4%)
1	F	0.29	0/539	0.70	0/730
1	G	0.28	0/539	0.71	0/730
1	H	0.29	0/539	0.70	0/730
1	I	0.29	0/539	0.75	0/730
All	All	0.30	0/4870	0.75	3/6595 (0.0%)

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	E	22	ALA	CA-C-N	-6.60	111.64	122.54
1	E	22	ALA	C-N-CA	-6.60	111.64	122.54
1	E	66	GLU	CB-CA-C	-5.55	99.06	109.72

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	525	0	514	8	0
1	B	544	0	531	4	0
1	C	525	0	514	12	0
1	D	525	0	514	3	0
1	E	525	0	514	13	0
1	F	525	0	514	5	0
1	G	525	0	514	9	0
1	H	525	0	514	8	0
1	I	525	0	514	5	0
2	J	54	0	19	0	0
2	K	54	0	20	0	0
2	L	54	0	20	1	0
3	A	11	0	10	0	0
3	C	11	0	10	0	0
3	D	11	0	10	0	0
3	F	11	0	10	0	0
3	H	11	0	10	0	0
3	I	11	0	10	0	0
4	A	12	0	11	0	0
4	C	12	0	11	1	0
4	E	12	0	11	0	0
4	G	12	0	11	0	0
4	H	12	0	11	1	0
5	A	5	0	0	0	0
5	E	5	0	0	0	0
5	F	5	0	0	0	0
5	G	5	0	0	0	0
5	H	5	0	0	0	0
6	A	2	0	0	0	0
6	B	2	0	0	0	0
6	C	4	0	0	1	0
6	D	2	0	0	0	0
6	E	1	0	0	0	0
6	F	3	0	0	0	0
6	G	2	0	0	0	0
6	H	2	0	0	0	0
6	I	2	0	0	0	0
7	B	8	0	12	0	0
8	A	24	0	0	0	0
8	B	20	0	0	0	0
8	C	21	0	0	0	0
8	D	19	0	0	0	0
8	E	32	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
8	F	19	0	0	0	0
8	G	28	0	0	0	0
8	H	24	0	0	0	0
8	I	23	0	0	0	0
All	All	5295	0	4829	48	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 5.

All (48) 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:H:45:LYS:NZ	1:H:47:ARG:HD2	1.97	0.79
1:C:66:GLU:HB2	1:E:23:HIS:HD2	1.54	0.72
1:H:45:LYS:HZ1	1:H:47:ARG:HD2	1.58	0.66
1:A:23:HIS:HA	1:A:44:ARG:HD3	1.78	0.65
1:E:59:ARG:NH1	1:G:29:TYR:OH	2.33	0.61
1:B:31:SER:OG	1:B:33:LYS:HE2	2.02	0.60
2:L:3[A]:SGN:O4	2:L:3[A]:SGN:O6	2.20	0.58
1:C:66:GLU:HB2	1:E:23:HIS:CD2	2.37	0.57
1:A:67:MET:HE2	1:C:23:HIS:HD2	1.70	0.56
1:C:63:ASN:HA	1:E:23:HIS:HE2	1.69	0.56
1:B:23:HIS:HD2	1:D:67:MET:HE2	1.72	0.55
1:C:66:GLU:HA	1:E:45:LYS:HG2	1.90	0.54
1:E:6:ASP:OD1	1:F:47:ARG:NE	2.40	0.53
1:F:24:ILE:HD12	1:F:41:PHE:HB3	1.90	0.52
1:H:45:LYS:HZ2	1:H:47:ARG:HD2	1.69	0.52
1:G:23:HIS:HD2	1:I:67:MET:HE2	1.75	0.52
1:A:66:GLU:O	1:C:44:ARG:NH1	2.43	0.51
1:H:52:ASN:HB3	1:H:55:LYS:HG2	1.92	0.51
1:A:54:GLU:O	1:A:59:ARG:NH1	2.44	0.50
1:C:66:GLU:CB	1:E:23:HIS:HD2	2.22	0.50
1:F:26:GLU:HB3	1:F:42:VAL:HB	1.95	0.48
1:A:24:ILE:HD12	1:A:41:PHE:HB3	1.95	0.48
1:E:20:PRO:HB2	1:E:23:HIS:CE1	2.48	0.48
1:G:63:ASN:O	1:G:67:MET:HG2	2.14	0.47
1:G:39:VAL:HG23	1:G:53:PRO:HG3	1.97	0.46
1:H:44:ARG:HE	1:H:44:ARG:HB2	1.51	0.45
1:H:20:PRO:HG2	1:H:23:HIS:CG	2.52	0.45
1:B:23:HIS:CD2	1:D:67:MET:HE2	2.52	0.44
1:G:13:ALA:HB2	1:I:33:LYS:HD2	2.00	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:33:LYS:HG3	4:C:102:BGC:O2	2.19	0.42
1:C:33:LYS:HD2	1:E:13:ALA:HB2	2.01	0.42
1:F:56:LYS:HE3	1:F:60:GLU:OE2	2.20	0.42
1:G:47:ARG:NH2	1:I:26:GLU:HB2	2.34	0.42
1:E:27:TYR:CZ	1:E:62:ILE:HG23	2.55	0.42
1:G:26:GLU:HG2	1:G:42:VAL:HB	2.01	0.42
1:A:54:GLU:HG2	1:C:17:ARG:HH12	1.85	0.42
4:H:102:BGC:O4	4:H:102:BGC:O6	2.36	0.41
1:G:47:ARG:HH21	1:I:26:GLU:HB2	1.85	0.41
1:E:23:HIS:N	1:E:23:HIS:ND1	2.67	0.41
1:H:43:THR:OG1	1:H:45:LYS:HG2	2.21	0.41
1:B:27:TYR:CZ	1:B:62:ILE:HG23	2.55	0.41
1:A:52:ASN:HB3	1:A:55:LYS:HG2	2.01	0.41
1:C:66:GLU:HG3	1:E:23:HIS:CD2	2.56	0.41
6:C:106:CL:CL	1:E:20:PRO:HG2	2.58	0.41
1:A:67:MET:HE2	1:C:23:HIS:CD2	2.53	0.40
1:G:23:HIS:CD2	1:I:67:MET:HE2	2.54	0.40
1:H:8:THR:HA	1:H:9:PRO:HD3	1.96	0.40
1:D:23:HIS:CD2	1:F:67:MET:HE2	2.56	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	63/65 (97%)	63 (100%)	0	0	100	100
1	B	65/65 (100%)	65 (100%)	0	0	100	100
1	C	63/65 (97%)	62 (98%)	1 (2%)	0	100	100
1	D	63/65 (97%)	63 (100%)	0	0	100	100
1	E	63/65 (97%)	63 (100%)	0	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	F	63/65 (97%)	63 (100%)	0	0	100	100
1	G	63/65 (97%)	63 (100%)	0	0	100	100
1	H	63/65 (97%)	63 (100%)	0	0	100	100
1	I	63/65 (97%)	63 (100%)	0	0	100	100
All	All	569/585 (97%)	568 (100%)	1 (0%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains [i](#)

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	59/59 (100%)	59 (100%)	0	100	100
1	B	61/59 (103%)	61 (100%)	0	100	100
1	C	59/59 (100%)	57 (97%)	2 (3%)	32	48
1	D	59/59 (100%)	58 (98%)	1 (2%)	53	72
1	E	59/59 (100%)	58 (98%)	1 (2%)	53	72
1	F	59/59 (100%)	59 (100%)	0	100	100
1	G	59/59 (100%)	59 (100%)	0	100	100
1	H	59/59 (100%)	59 (100%)	0	100	100
1	I	59/59 (100%)	58 (98%)	1 (2%)	53	72
All	All	533/531 (100%)	528 (99%)	5 (1%)	70	82

All (5) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	C	4	SER
1	C	46	ASN
1	D	46	ASN
1	E	23	HIS
1	I	46	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	36	ASN
1	E	23	HIS
1	E	63	ASN
1	F	23	HIS
1	F	63	ASN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

9 monosaccharides are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	SGN	J	1[A]	2	19,19,20	4.51	2 (10%)	23,29,31	3.40	11 (47%)
2	IDS	J	2[A]	2	16,16,17	1.48	2 (12%)	16,24,26	3.05	5 (31%)
2	SGN	J	3[A]	2	19,19,20	4.58	2 (10%)	23,29,31	3.08	7 (30%)
2	SGN	K	1[A]	2	19,19,20	4.56	2 (10%)	23,29,31	3.33	10 (43%)
2	IDS	K	2[A]	2	16,16,17	1.20	2 (12%)	16,24,26	3.24	5 (31%)
2	SGN	K	3[A]	2	19,19,20	4.59	2 (10%)	23,29,31	3.09	7 (30%)
2	SGN	L	1[A]	2	19,19,20	1.54	3 (15%)	23,29,31	3.24	9 (39%)
2	IDS	L	2[A]	2	16,16,17	1.16	2 (12%)	16,24,26	3.26	5 (31%)
2	SGN	L	3[A]	2	19,19,20	1.22	2 (10%)	23,29,31	3.15	8 (34%)

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	SGN	J	1[A]	2	-	6/11/28/31	0/1/1/1
2	IDS	J	2[A]	2	-	2/9/26/29	0/1/1/1
2	SGN	J	3[A]	2	-	2/11/28/31	0/1/1/1
2	SGN	K	1[A]	2	-	7/11/28/31	0/1/1/1
2	IDS	K	2[A]	2	-	1/9/26/29	0/1/1/1
2	SGN	K	3[A]	2	-	2/11/28/31	0/1/1/1
2	SGN	L	1[A]	2	-	5/11/28/31	0/1/1/1
2	IDS	L	2[A]	2	-	1/9/26/29	0/1/1/1
2	SGN	L	3[A]	2	-	9/11/28/31	0/1/1/1

All (19) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	K	1[A]	SGN	O1S-S1	19.53	1.64	1.42
2	J	1[A]	SGN	O2S-S1	19.28	1.64	1.42
2	K	3[A]	SGN	O2S-S1	19.27	1.64	1.42
2	J	3[A]	SGN	O1S-S1	19.25	1.64	1.42
2	L	1[A]	SGN	O5S-S2	4.31	1.64	1.45
2	K	3[A]	SGN	O4S-S2	4.31	1.64	1.45
2	J	3[A]	SGN	O5S-S2	4.27	1.63	1.45
2	J	2[A]	IDS	O2S-S	4.18	1.63	1.45
2	L	3[A]	SGN	O3S-S1	3.43	1.64	1.47
2	L	1[A]	SGN	O3S-S1	3.42	1.64	1.47
2	K	2[A]	IDS	O2-C2	-2.80	1.43	1.47
2	J	2[A]	IDS	O2-C2	-2.73	1.43	1.47
2	L	2[A]	IDS	O2-C2	-2.69	1.43	1.47
2	L	3[A]	SGN	O6S-S2	2.17	1.64	1.50
2	K	1[A]	SGN	O6S-S2	2.17	1.64	1.50
2	J	1[A]	SGN	O6S-S2	2.16	1.63	1.50
2	L	2[A]	IDS	O3S-S	2.15	1.63	1.50
2	K	2[A]	IDS	O3S-S	2.14	1.63	1.50
2	L	1[A]	SGN	O2S-S1	2.04	1.44	1.42

All (67) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	K	1[A]	SGN	O6S-S2-O6	-8.38	87.12	106.37

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	L	3[A]	SGN	O6S-S2-O6	-8.25	87.42	106.37
2	L	2[A]	IDS	O3S-S-O2	-8.23	87.47	106.37
2	J	1[A]	SGN	O6S-S2-O6	-8.21	87.50	106.37
2	K	2[A]	IDS	O3S-S-O2	-8.18	87.58	106.37
2	J	2[A]	IDS	O2S-S-O1S	-6.70	86.40	112.24
2	K	3[A]	SGN	O6-S2-O4S	-6.67	86.55	106.92
2	L	1[A]	SGN	O6-S2-O5S	-6.63	86.68	106.92
2	K	1[A]	SGN	O6S-S2-O5S	-6.62	85.39	108.56
2	J	3[A]	SGN	O5S-S2-O4S	-6.61	86.75	112.24
2	L	1[A]	SGN	O5S-S2-O4S	-6.55	86.96	112.24
2	K	3[A]	SGN	O5S-S2-O4S	-6.55	86.98	112.24
2	J	3[A]	SGN	O6-S2-O5S	-6.52	87.01	106.92
2	J	2[A]	IDS	O3S-S-O2S	-6.43	86.06	108.56
2	K	2[A]	IDS	O3S-S-O1S	-6.43	86.07	108.56
2	L	2[A]	IDS	O3S-S-O1S	-6.39	86.18	108.56
2	L	1[A]	SGN	O6S-S2-O5S	-6.33	86.42	108.56
2	K	1[A]	SGN	O6S-S2-O4S	-6.31	86.49	108.56
2	K	3[A]	SGN	O6S-S2-O4S	-6.30	86.50	108.56
2	L	3[A]	SGN	O6S-S2-O4S	-6.29	86.56	108.56
2	J	1[A]	SGN	O6S-S2-O4S	-6.25	86.67	108.56
2	L	3[A]	SGN	O6S-S2-O5S	-6.08	87.26	108.56
2	J	1[A]	SGN	O6S-S2-O5S	-6.04	87.42	108.56
2	J	3[A]	SGN	O6S-S2-O5S	-6.01	87.54	108.56
2	K	2[A]	IDS	O3S-S-O2S	-6.00	87.55	108.56
2	L	2[A]	IDS	O3S-S-O2S	-5.99	87.59	108.56
2	L	1[A]	SGN	O6S-S2-O6	5.97	120.10	106.37
2	K	3[A]	SGN	O6S-S2-O6	5.93	119.99	106.37
2	J	3[A]	SGN	O6S-S2-O6	5.77	119.64	106.37
2	J	2[A]	IDS	O3S-S-O2	5.63	119.30	106.37
2	J	1[A]	SGN	O1S-S1-O2S	-5.37	108.42	120.36
2	J	3[A]	SGN	O1S-S1-O2S	-5.01	109.22	120.36
2	L	3[A]	SGN	O1S-S1-O2S	-4.93	109.39	120.36
2	K	1[A]	SGN	O1S-S1-O2S	-4.92	109.41	120.36
2	K	3[A]	SGN	O1S-S1-O2S	-4.87	109.53	120.36
2	L	1[A]	SGN	O1S-S1-O2S	-4.86	109.54	120.36
2	L	3[A]	SGN	O6-S2-O5S	4.36	120.24	106.92
2	J	1[A]	SGN	O6-S2-O5S	4.22	119.80	106.92
2	J	3[A]	SGN	O6-S2-O4S	4.21	119.79	106.92
2	J	1[A]	SGN	O6-S2-O4S	4.17	119.66	106.92
2	K	1[A]	SGN	O6-S2-O5S	4.13	119.54	106.92
2	K	3[A]	SGN	O6-S2-O5S	4.08	119.37	106.92
2	K	1[A]	SGN	O6-S2-O4S	4.03	119.22	106.92

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	L	1[A]	SGN	O6-S2-O4S	4.02	119.20	106.92
2	L	3[A]	SGN	O6-S2-O4S	4.00	119.12	106.92
2	J	1[A]	SGN	C3-C2-N2	-3.55	105.66	110.32
2	J	1[A]	SGN	O4-C4-C3	-3.32	102.54	110.38
2	J	2[A]	IDS	O3S-S-O1S	3.32	120.17	108.56
2	L	2[A]	IDS	O2-C2-C3	3.28	111.54	106.95
2	J	1[A]	SGN	C1-O5-C5	3.27	116.57	112.19
2	J	2[A]	IDS	O2-C2-C3	3.24	111.47	106.95
2	J	3[A]	SGN	O6S-S2-O4S	3.22	119.81	108.56
2	L	1[A]	SGN	O6S-S2-O4S	3.19	119.71	108.56
2	K	3[A]	SGN	O6S-S2-O5S	3.17	119.64	108.56
2	K	2[A]	IDS	O2-C2-C3	3.13	111.33	106.95
2	L	3[A]	SGN	C3-C2-N2	-2.95	106.45	110.32
2	K	1[A]	SGN	C1-O5-C5	2.92	116.10	112.19
2	J	1[A]	SGN	C1-C2-N2	2.74	114.56	110.22
2	K	1[A]	SGN	O4-C4-C3	-2.74	103.91	110.38
2	L	3[A]	SGN	O5-C5-C6	2.72	113.47	107.59
2	L	1[A]	SGN	C1-O5-C5	2.72	115.83	112.19
2	L	1[A]	SGN	O4-C4-C3	-2.49	104.51	110.38
2	K	1[A]	SGN	O5-C5-C6	-2.34	102.54	107.59
2	J	1[A]	SGN	O6-C6-C5	2.21	111.50	107.57
2	K	2[A]	IDS	O2S-S-O1S	2.08	120.28	112.24
2	K	1[A]	SGN	O5S-S2-O4S	2.02	120.02	112.24
2	L	2[A]	IDS	O2S-S-O1S	2.01	119.97	112.24

There are no chirality outliers.

All (35) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	J	1[A]	SGN	C1-C2-N2-S1
2	J	1[A]	SGN	C4-C5-C6-O6
2	J	1[A]	SGN	O5-C5-C6-O6
2	J	3[A]	SGN	C6-O6-S2-O6S
2	K	1[A]	SGN	C1-C2-N2-S1
2	K	1[A]	SGN	C4-C5-C6-O6
2	L	1[A]	SGN	C1-C2-N2-S1
2	L	2[A]	IDS	C2-O2-S-O1S
2	L	3[A]	SGN	C4-C5-C6-O6
2	L	3[A]	SGN	O5-C5-C6-O6
2	L	3[A]	SGN	C2-N2-S1-O1S
2	L	3[A]	SGN	C2-N2-S1-O2S
2	L	3[A]	SGN	C2-N2-S1-O3S

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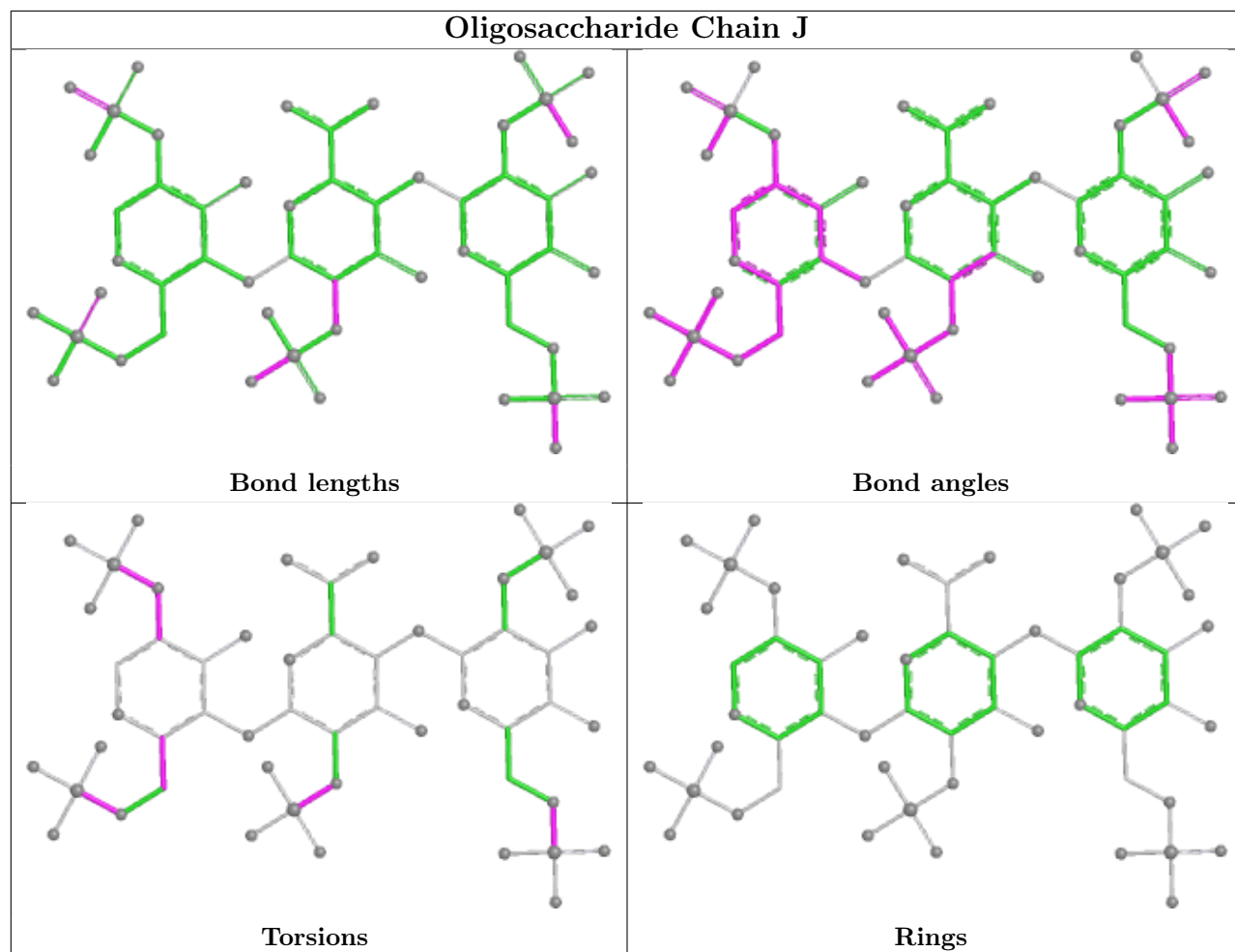
Mol	Chain	Res	Type	Atoms
2	L	3[A]	SGN	C6-O6-S2-O4S
2	J	1[A]	SGN	C6-O6-S2-O4S
2	J	1[A]	SGN	C6-O6-S2-O5S
2	K	3[A]	SGN	C6-O6-S2-O4S
2	J	2[A]	IDS	C2-O2-S-O1S
2	J	2[A]	IDS	C2-O2-S-O2S
2	K	2[A]	IDS	C2-O2-S-O1S
2	K	3[A]	SGN	C6-O6-S2-O6S
2	L	1[A]	SGN	C6-O6-S2-O6S
2	L	3[A]	SGN	C1-C2-N2-S1
2	K	1[A]	SGN	C6-O6-S2-O5S
2	L	1[A]	SGN	C6-O6-S2-O5S
2	J	3[A]	SGN	C6-O6-S2-O5S
2	J	1[A]	SGN	C2-N2-S1-O2S
2	L	3[A]	SGN	C6-O6-S2-O6S
2	K	1[A]	SGN	C6-O6-S2-O4S
2	L	3[A]	SGN	C3-C2-N2-S1
2	K	1[A]	SGN	O5-C5-C6-O6
2	L	1[A]	SGN	C4-C5-C6-O6
2	K	1[A]	SGN	C5-C6-O6-S2
2	K	1[A]	SGN	C2-N2-S1-O1S
2	L	1[A]	SGN	C2-N2-S1-O2S

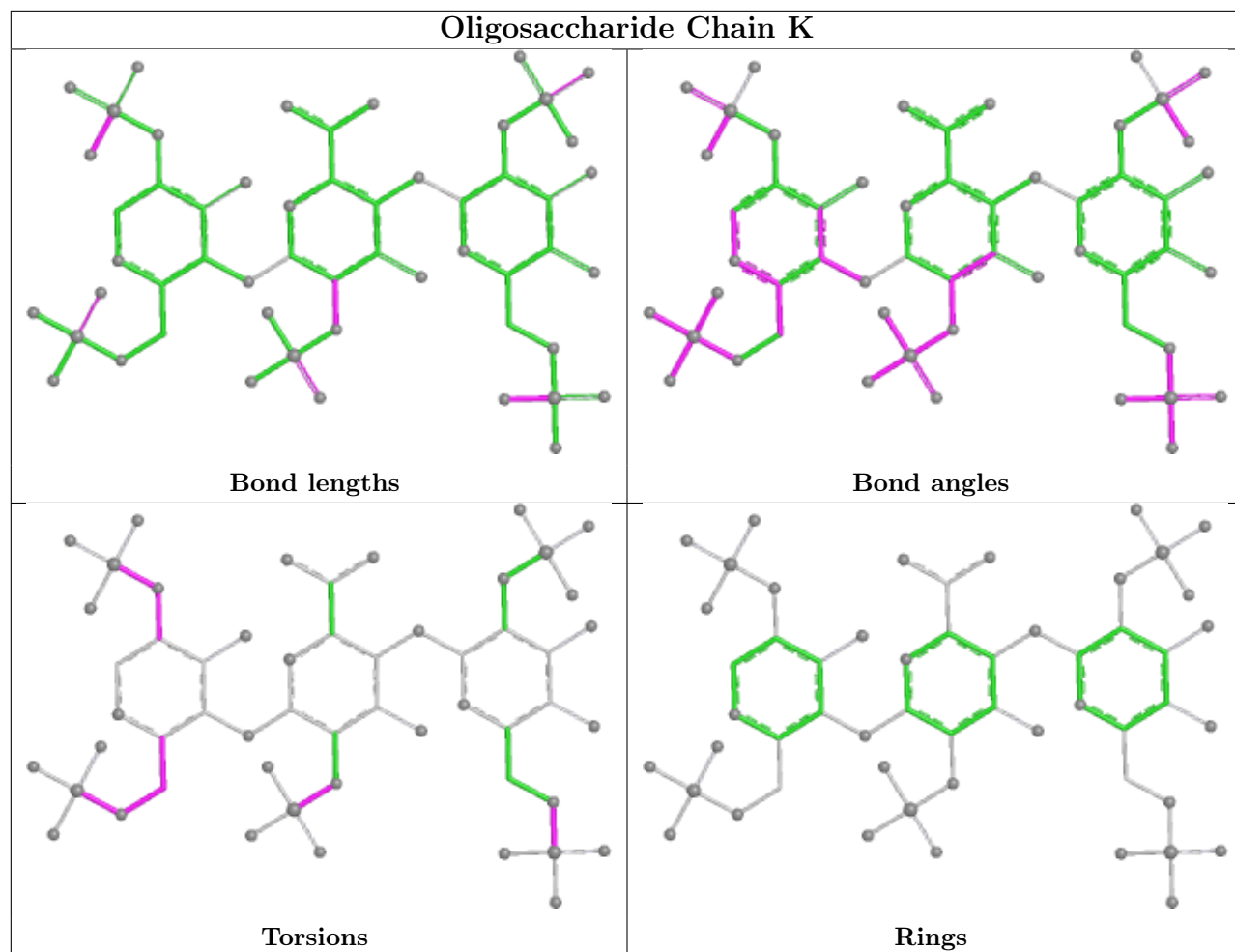
There are no ring outliers.

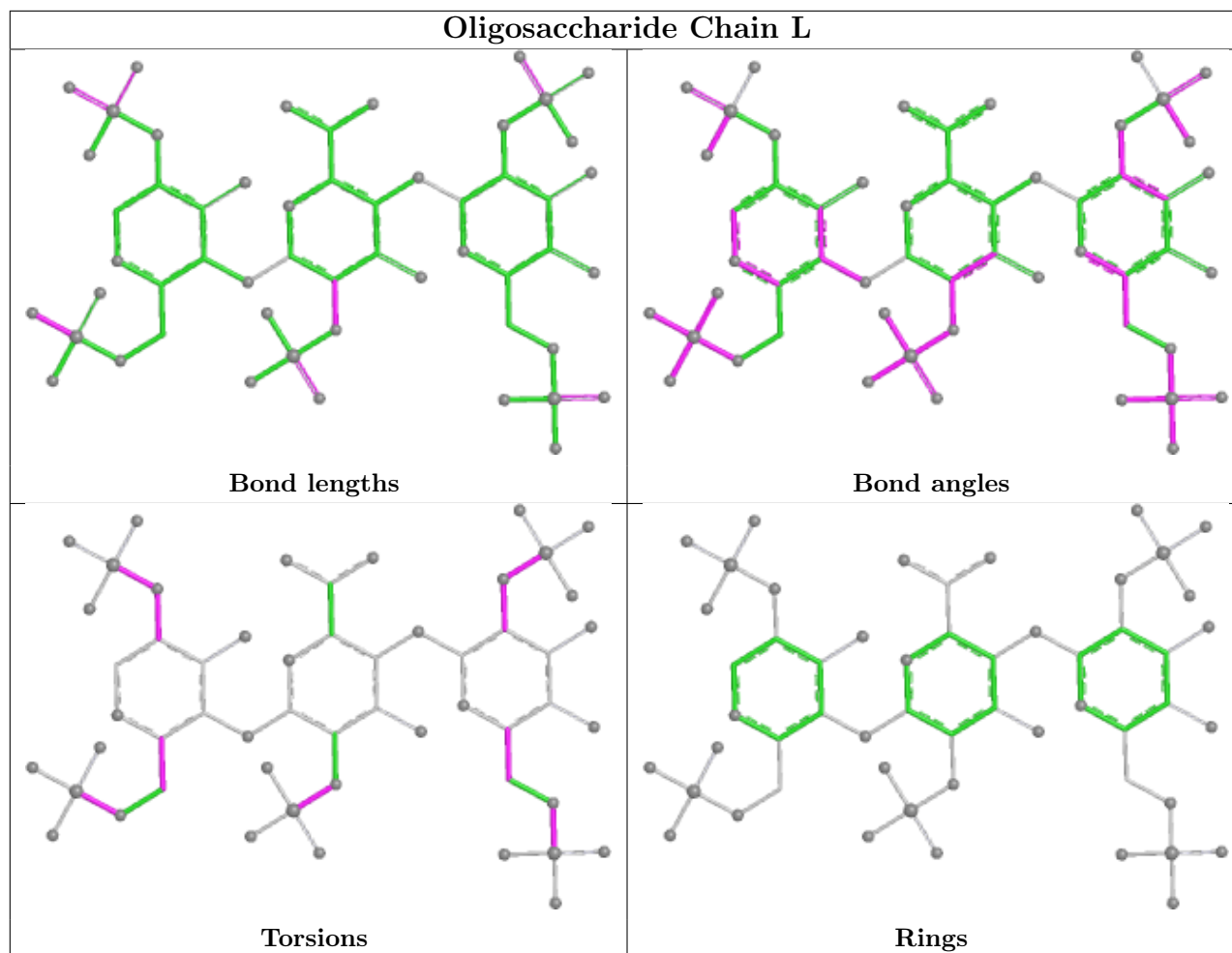
1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	L	3[A]	SGN	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 37 ligands modelled in this entry, 20 are monoatomic - leaving 17 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$
3	GLA	C	101	1	11,11,12	1.01	1 (9%)	15,15,17	1.40	3 (20%)
3	GLA	H	101	1	11,11,12	0.97	1 (9%)	15,15,17	1.12	2 (13%)
5	SO4	A	103	-	4,4,4	0.24	0	6,6,6	0.08	0
5	SO4	E	102	-	4,4,4	0.24	0	6,6,6	0.08	0
4	BGC	G	101	-	12,12,12	2.08	3 (25%)	17,17,17	0.93	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
5	SO4	F	102	-	4,4,4	0.23	0	6,6,6	0.09	0
5	SO4	H	103	-	4,4,4	0.23	0	6,6,6	0.08	0
4	BGC	A	102	-	12,12,12	2.10	3 (25%)	17,17,17	0.95	1 (5%)
4	BGC	C	102	-	12,12,12	2.11	5 (41%)	17,17,17	1.33	1 (5%)
3	GLA	D	101	1	11,11,12	0.90	0	15,15,17	1.22	1 (6%)
3	GLA	A	101	1	11,11,12	0.97	1 (9%)	15,15,17	1.18	2 (13%)
5	SO4	G	102	-	4,4,4	0.24	0	6,6,6	0.08	0
4	BGC	E	101	-	12,12,12	2.08	2 (16%)	17,17,17	1.02	1 (5%)
7	TRS	B	103	-	7,7,7	0.33	0	9,9,9	0.32	0
3	GLA	F	101	1	11,11,12	0.94	0	15,15,17	1.18	2 (13%)
3	GLA	I	101	1	11,11,12	0.91	0	15,15,17	1.22	1 (6%)
4	BGC	H	102	-	12,12,12	2.14	4 (33%)	17,17,17	0.79	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
3	GLA	C	101	1	-	0/2/19/22	0/1/1/1
3	GLA	H	101	1	-	2/2/19/22	0/1/1/1
4	BGC	G	101	-	-	2/2/22/22	0/1/1/1
4	BGC	A	102	-	-	2/2/22/22	0/1/1/1
4	BGC	C	102	-	-	2/2/22/22	0/1/1/1
3	GLA	D	101	1	-	0/2/19/22	0/1/1/1
3	GLA	A	101	1	-	2/2/19/22	0/1/1/1
4	BGC	E	101	-	-	1/2/22/22	0/1/1/1
7	TRS	B	103	-	-	1/9/9/9	-
3	GLA	F	101	1	-	2/2/19/22	0/1/1/1
3	GLA	I	101	1	-	0/2/19/22	0/1/1/1
4	BGC	H	102	-	-	2/2/22/22	0/1/1/1

All (20) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	H	102	BGC	O5-C1	4.85	1.54	1.42
4	E	101	BGC	O5-C1	4.83	1.54	1.42
4	G	101	BGC	O5-C1	4.81	1.54	1.42
4	A	102	BGC	O5-C1	4.80	1.54	1.42

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	C	102	BGC	O5-C1	4.05	1.52	1.42
4	C	102	BGC	O1-C1	-3.57	1.28	1.39
4	A	102	BGC	O1-C1	-2.94	1.30	1.39
4	H	102	BGC	O1-C1	-2.91	1.30	1.39
4	E	101	BGC	O1-C1	-2.89	1.30	1.39
4	G	101	BGC	O1-C1	-2.86	1.30	1.39
4	C	102	BGC	C4-C5	-2.82	1.47	1.53
4	C	102	BGC	C4-C3	-2.63	1.45	1.52
4	H	102	BGC	C4-C3	-2.22	1.46	1.52
4	C	102	BGC	C3-C2	-2.15	1.46	1.52
3	C	101	GLA	C2-C3	2.14	1.55	1.52
4	A	102	BGC	C4-C3	-2.08	1.46	1.52
4	G	101	BGC	C4-C3	-2.05	1.47	1.52
3	H	101	GLA	C2-C3	2.05	1.55	1.52
4	H	102	BGC	C4-C5	-2.04	1.48	1.53
3	A	101	GLA	C2-C3	2.03	1.55	1.52

All (14) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	D	101	GLA	C1-O5-C5	3.90	117.41	112.19
3	I	101	GLA	C1-O5-C5	3.80	117.27	112.19
4	C	102	BGC	O3-C3-C2	-3.37	102.43	110.38
3	A	101	GLA	C1-O5-C5	3.12	116.37	112.19
3	F	101	GLA	C1-O5-C5	3.12	116.36	112.19
3	H	101	GLA	C1-O5-C5	2.87	116.03	112.19
3	C	101	GLA	C2-C3-C4	2.68	115.57	110.86
3	C	101	GLA	C1-O5-C5	2.46	115.48	112.19
3	C	101	GLA	C3-C4-C5	2.35	114.49	110.23
3	A	101	GLA	O2-C2-C1	2.17	114.19	109.22
3	F	101	GLA	O2-C2-C1	2.14	114.13	109.22
4	A	102	BGC	O5-C1-C2	-2.06	106.68	110.30
4	E	101	BGC	O5-C1-C2	-2.04	106.71	110.30
3	H	101	GLA	O2-C2-C1	2.03	113.86	109.22

There are no chirality outliers.

All (16) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	C	102	BGC	C4-C5-C6-O6
4	C	102	BGC	O5-C5-C6-O6
4	H	102	BGC	C4-C5-C6-O6

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Mol	Chain	Res	Type	Atoms
4	A	102	BGC	O5-C5-C6-O6
4	H	102	BGC	O5-C5-C6-O6
4	A	102	BGC	C4-C5-C6-O6
4	G	101	BGC	O5-C5-C6-O6
3	H	101	GLA	O5-C5-C6-O6
3	A	101	GLA	O5-C5-C6-O6
3	F	101	GLA	O5-C5-C6-O6
4	E	101	BGC	O5-C5-C6-O6
3	H	101	GLA	C4-C5-C6-O6
3	A	101	GLA	C4-C5-C6-O6
4	G	101	BGC	C4-C5-C6-O6
3	F	101	GLA	C4-C5-C6-O6
7	B	103	TRS	C3-C-C1-O1

There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	C	102	BGC	1	0
4	H	102	BGC	1	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data [i](#)

6.1 Protein, DNA and RNA chains [i](#)

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	65/65 (100%)	-1.19	0 100 100	25, 36, 63, 77	0
1	B	65/65 (100%)	-1.27	0 100 100	17, 34, 52, 74	3 (4%)
1	C	65/65 (100%)	-1.30	0 100 100	21, 35, 55, 73	0
1	D	65/65 (100%)	-1.31	0 100 100	22, 37, 54, 71	0
1	E	65/65 (100%)	-1.24	0 100 100	24, 34, 57, 71	1 (1%)
1	F	65/65 (100%)	-1.19	0 100 100	24, 37, 63, 76	0
1	G	65/65 (100%)	-1.30	0 100 100	23, 34, 54, 77	1 (1%)
1	H	65/65 (100%)	-1.19	0 100 100	24, 36, 67, 75	0
1	I	65/65 (100%)	-1.31	0 100 100	21, 36, 54, 69	0
All	All	585/585 (100%)	-1.25	0 100 100	17, 36, 61, 77	5 (0%)

There are no RSRZ outliers to report.

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.

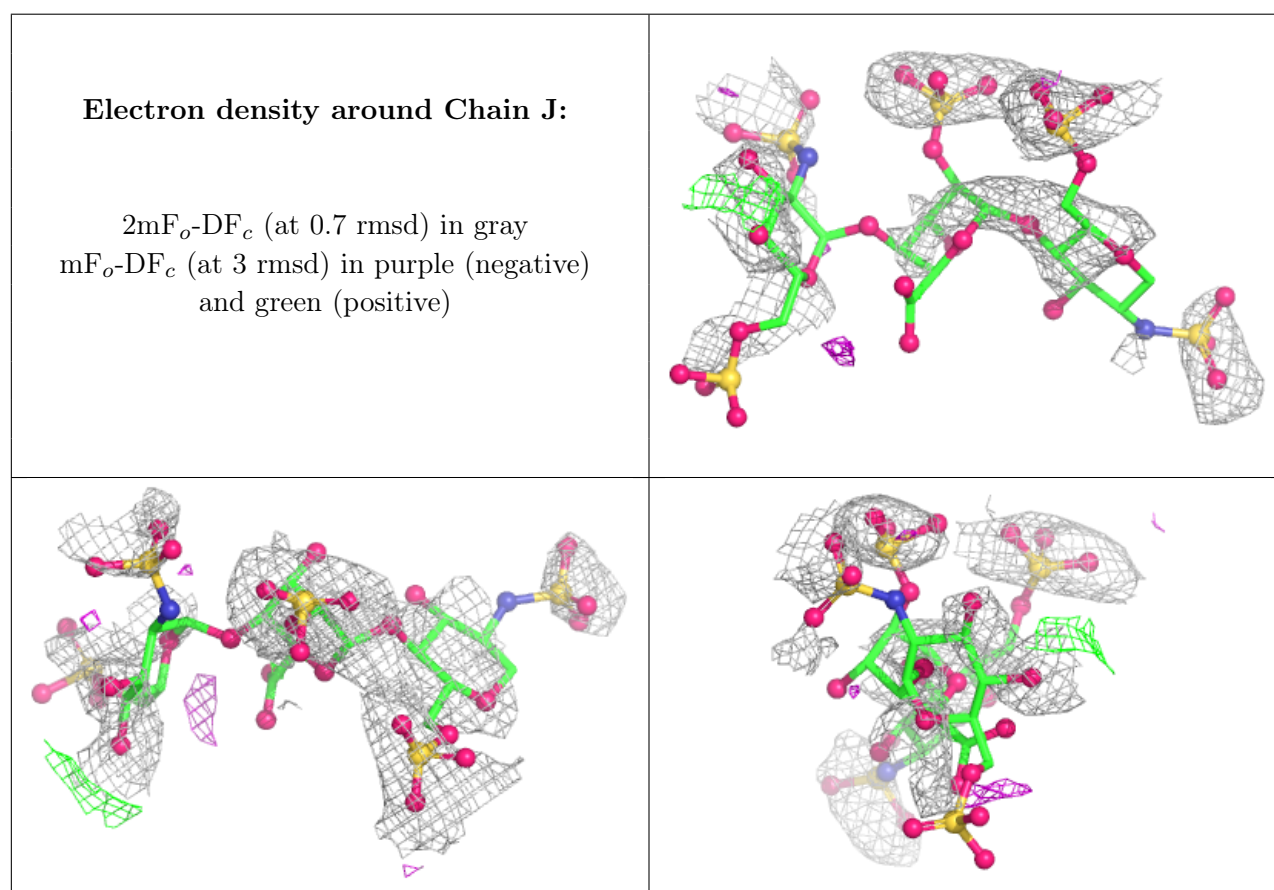
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	SGN	L	3[A]	19/20	0.90	0.17	101,117,131,141	19
2	SGN	K	3[A]	19/20	0.91	0.14	97,124,151,156	19

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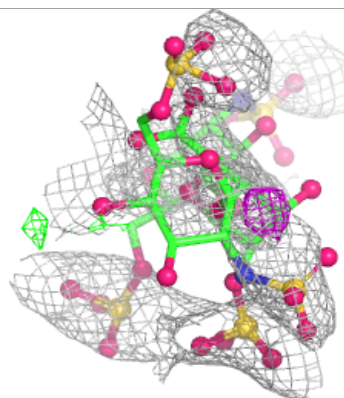
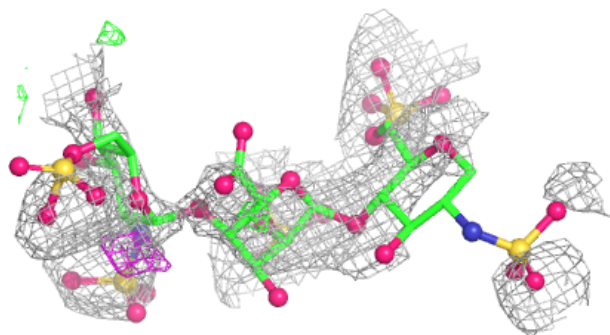
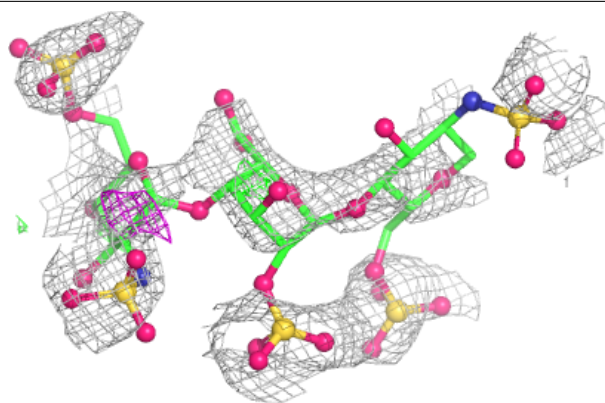
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
2	SGN	J	3[A]	19/20	0.92	0.15	97,120,147,152	19
2	IDS	L	2[A]	16/17	0.93	0.17	76,113,120,122	16
2	IDS	J	2[A]	16/17	0.95	0.17	76,116,120,122	16
2	SGN	K	1[A]	19/20	0.96	0.10	107,124,145,166	0
2	SGN	J	1[A]	19/20	0.96	0.11	101,110,141,144	0
2	IDS	K	2[A]	16/17	0.97	0.14	70,120,123,125	16
2	SGN	L	1[A]	19/20	0.97	0.12	98,117,129,150	19

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.

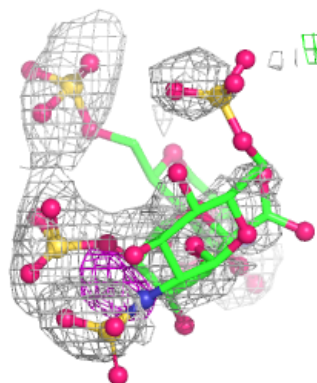
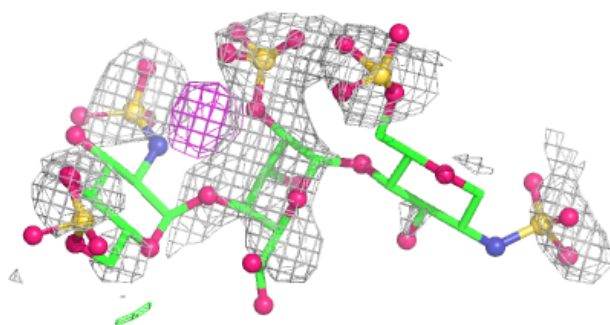
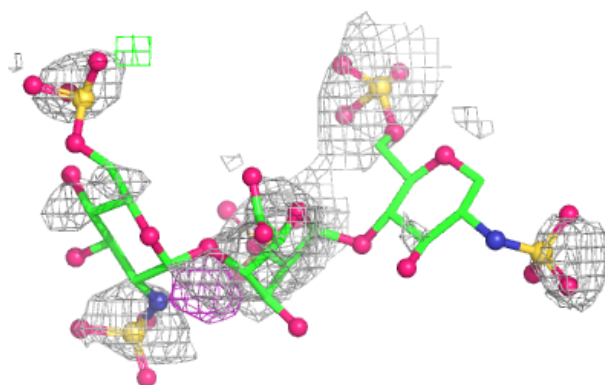


Electron density around Chain K:

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

**Electron density around Chain L:**

$2mF_o-DF_c$ (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	SO4	A	103	5/5	0.96	0.10	58,65,90,95	5
5	SO4	E	102	5/5	0.96	0.11	71,89,145,160	0
5	SO4	G	102	5/5	0.96	0.09	67,125,155,164	0
6	CL	C	105	1/1	0.96	0.16	73,73,73,73	0
6	CL	D	103	1/1	0.96	0.12	74,74,74,74	1
6	CL	G	103	1/1	0.96	0.09	80,80,80,80	0
6	CL	E	103	1/1	0.97	0.09	78,78,78,78	0
6	CL	C	104	1/1	0.97	0.10	81,81,81,81	0
6	CL	I	102	1/1	0.97	0.09	79,79,79,79	0
6	CL	I	103	1/1	0.97	0.08	63,63,63,63	1
7	TRS	B	103	8/8	0.97	0.07	57,74,90,99	0
6	CL	C	106	1/1	0.98	0.12	50,50,50,50	1
6	CL	D	102	1/1	0.98	0.07	82,82,82,82	0
5	SO4	H	103	5/5	0.98	0.06	54,59,87,91	5
6	CL	A	105	1/1	0.98	0.07	30,30,30,30	0
6	CL	F	105	1/1	0.98	0.09	74,74,74,74	0
6	CL	B	101	1/1	0.98	0.08	81,81,81,81	0
6	CL	G	104	1/1	0.98	0.07	83,83,83,83	0
6	CL	H	104	1/1	0.98	0.06	67,67,67,67	0
6	CL	B	102	1/1	0.98	0.07	76,76,76,76	0
5	SO4	F	102	5/5	0.98	0.07	51,53,84,90	5
4	BGC	E	101	12/12	0.98	0.06	53,76,89,89	0
4	BGC	H	102	12/12	0.99	0.05	50,59,68,69	0
3	GLA	C	101	11/12	0.99	0.06	40,48,62,98	0
3	GLA	D	101	11/12	0.99	0.06	41,53,74,107	0
3	GLA	F	101	11/12	0.99	0.06	44,61,75,90	0
3	GLA	H	101	11/12	0.99	0.06	45,63,79,90	0
6	CL	F	103	1/1	0.99	0.12	44,44,44,44	0
6	CL	F	104	1/1	0.99	0.04	34,34,34,34	1
3	GLA	I	101	11/12	0.99	0.05	40,45,69,102	0
6	CL	A	104	1/1	0.99	0.06	32,32,32,32	1
4	BGC	A	102	12/12	0.99	0.05	45,62,73,73	0
4	BGC	C	102	12/12	0.99	0.07	47,64,80,93	0
6	CL	H	105	1/1	0.99	0.06	25,25,25,25	1
3	GLA	A	101	11/12	0.99	0.05	43,64,75,89	0
6	CL	C	103	1/1	0.99	0.13	53,53,53,53	0
4	BGC	G	101	12/12	0.99	0.05	63,80,92,95	0

6.5 Other polymers [i](#)

There are no such residues in this entry.