



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 09:22 PM UTC

PDB ID : 3M2M / pdb_00003m2m
Title : Rat galectin-1 complex with lactose
Authors : Lobsanov, Y.D.; Rini, J.M.; Leffler, H.
Deposited on : 2010-03-07
Resolution : 2.95 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

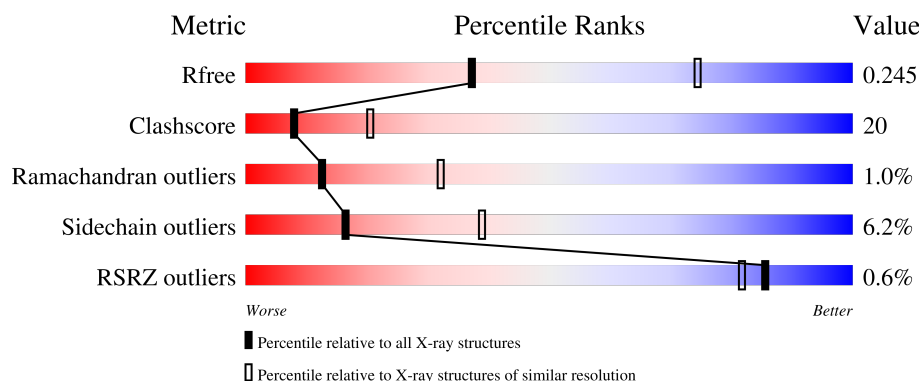
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.95 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	1130 (2.98-2.94)
Clashscore	190562	1157 (2.98-2.94)
Ramachandran outliers	187476	1101 (2.98-2.94)
Sidechain outliers	187428	1101 (2.98-2.94)
RSRZ outliers	180081	1130 (2.98-2.94)

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	134	% 60% 35% ..
1	B	134	55% 32% .. 9%
1	C	134	61% 34% ..
1	D	134	% 64% 31% ..
1	E	134	% 62% 33% ..

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Mol	Chain	Length	Quality of chain
1	F	134	% 57% 30% 9%
1	G	134	57% 31% 9%
1	H	134	60% 35%
2	I	2	50% 50%
2	J	2	100%
2	K	2	100%
2	L	2	100%
2	M	2	100%

2 Entry composition [i](#)

There are 2 unique types of molecules in this entry. The entry contains 8076 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 Galectin-1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	132	Total	C	N	O	S	0	0	0
			1024	643	175	199	7			
1	B	122	Total	C	N	O	S	0	0	0
			947	593	162	185	7			
1	C	132	Total	C	N	O	S	0	0	0
			1024	643	175	199	7			
1	D	132	Total	C	N	O	S	0	0	0
			1024	643	175	199	7			
1	E	132	Total	C	N	O	S	0	0	0
			1024	643	175	199	7			
1	F	122	Total	C	N	O	S	0	0	0
			947	593	162	185	7			
1	G	122	Total	C	N	O	S	0	0	0
			947	593	162	185	7			
1	H	132	Total	C	N	O	S	0	0	0
			1024	643	175	199	7			

- Molecule 2 is an oligosaccharide called beta-D-galactopyranose-(1-4)-beta-D-glucopyranose.



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace
2	I	2	Total	C	O	0	0	0
			23	12	11			
2	J	2	Total	C	O	0	0	0
			23	12	11			
2	K	2	Total	C	O	0	0	0
			23	12	11			
2	L	2	Total	C	O	0	0	0
			23	12	11			

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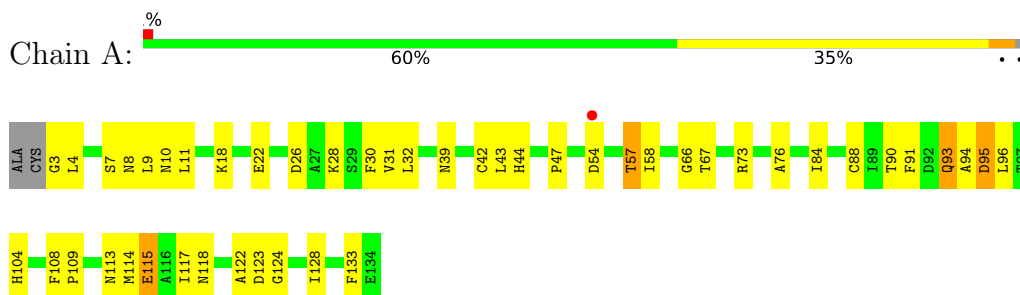
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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace
2	M	2	Total	C	O	0	0	0
			23	12	11			

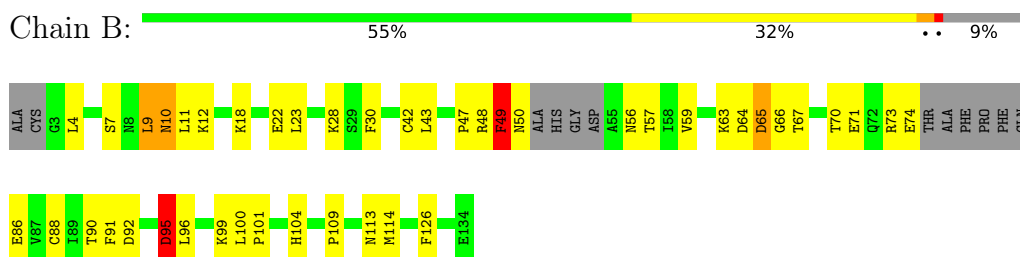
3 Residue-property plots [i](#)

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.

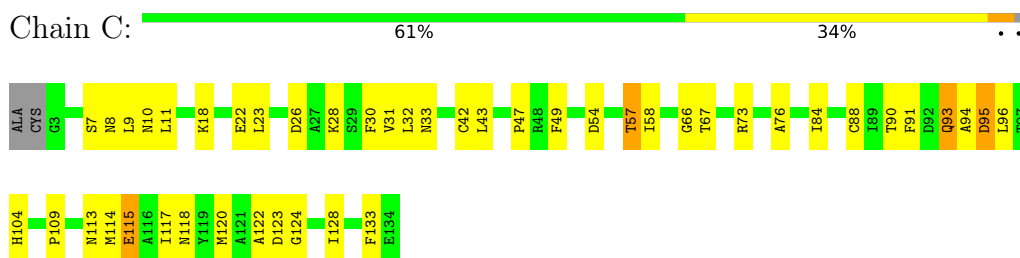
• Molecule 1: Galectin-1



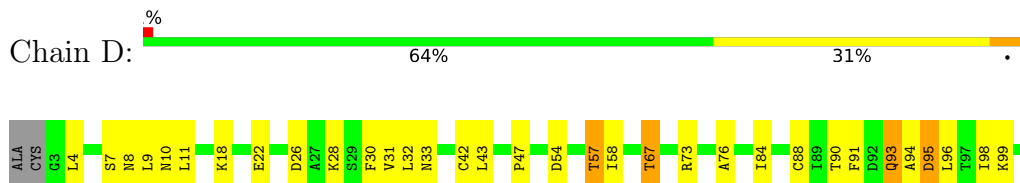
• Molecule 1: Galectin-1



• Molecule 1: Galectin-1

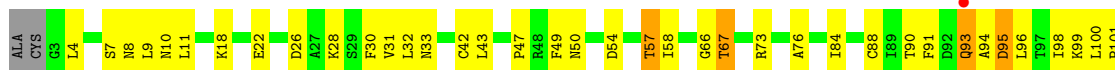


• Molecule 1: Galectin-1

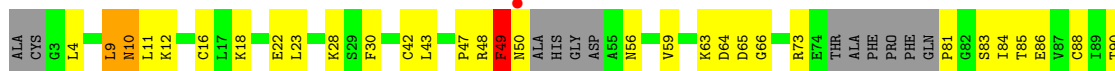




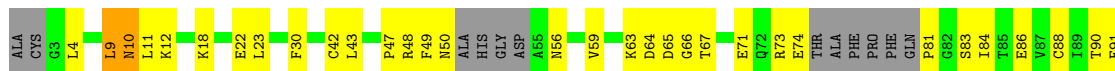
- Molecule 1: Galectin-1



- Molecule 1: Galectin-1



- Molecule 1: Galectin-1



- Molecule 1: Galectin-1



- Molecule 2: beta-D-galactopyranose-(1-4)-beta-D-glucopyranose





- Molecule 2: beta-D-galactopyranose-(1-4)-beta-D-glucopyranose

Chain J:  100%



- Molecule 2: beta-D-galactopyranose-(1-4)-beta-D-glucopyranose

Chain K:  100%



- Molecule 2: beta-D-galactopyranose-(1-4)-beta-D-glucopyranose

Chain L:  100%



- Molecule 2: beta-D-galactopyranose-(1-4)-beta-D-glucopyranose

Chain M:  100%



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	40.10Å 58.60Å 121.60Å 101.10° 91.60° 110.90°	Depositor
Resolution (Å)	36.00 – 2.95 36.00 – 2.95	Depositor EDS
% Data completeness (in resolution range)	73.7 (36.00-2.95) 73.7 (36.00-2.95)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.33 (at 2.81Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.206 , 0.248 0.203 , 0.245	Depositor DCC
R_{free} test set	1262 reflections (6.89%)	wwPDB-VP
Wilson B-factor (Å ²)	26.9	Xtriage
Anisotropy	0.541	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 30.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	0.052 for h,-h-k,-l	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	8076	wwPDB-VP
Average B, all atoms (Å ²)	27.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 29.30 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 1.5859e-03. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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: BGC, GAL, CSO

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	0.47	0/1030	0.95	2/1389 (0.1%)
1	B	0.49	0/947	0.98	2/1271 (0.2%)
1	C	0.47	0/1030	0.93	1/1389 (0.1%)
1	D	0.48	0/1030	0.94	0/1389
1	E	0.49	0/1030	0.95	1/1389 (0.1%)
1	F	0.48	0/947	0.99	2/1271 (0.2%)
1	G	0.51	0/947	0.92	1/1271 (0.1%)
1	H	0.47	0/1030	0.94	1/1389 (0.1%)
All	All	0.48	0/7991	0.95	10/10758 (0.1%)

There are no bond length outliers.

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	F	49	PHE	N-CA-C	-13.53	96.09	112.59
1	B	49	PHE	N-CA-C	-10.37	99.70	114.12
1	G	95	ASP	N-CA-C	6.35	117.87	108.60
1	B	95	ASP	N-CA-C	6.06	117.45	108.60
1	F	95	ASP	N-CA-C	5.83	117.11	108.60
1	C	49	PHE	N-CA-C	-5.74	105.11	111.36
1	H	49	PHE	N-CA-C	-5.16	105.73	111.36
1	A	108	PHE	CA-C-N	5.11	125.02	119.76
1	A	108	PHE	C-N-CA	5.11	125.02	119.76
1	E	49	PHE	N-CA-C	-5.02	105.89	111.36

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	1024	0	983	47	0
1	B	947	0	918	35	0
1	C	1024	0	983	44	0
1	D	1024	0	983	40	0
1	E	1024	0	983	47	0
1	F	947	0	918	36	0
1	G	947	0	918	31	0
1	H	1024	0	983	46	0
2	I	23	0	21	2	0
2	J	23	0	21	0	0
2	K	23	0	21	0	0
2	L	23	0	21	0	0
2	M	23	0	21	0	0
All	All	8076	0	7774	310	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 20.

All (310) 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:A:95:ASP:HB3	1:A:109:PRO:HA	1.53	0.91
1:D:95:ASP:HB3	1:D:109:PRO:HA	1.54	0.90
1:E:95:ASP:HB3	1:E:109:PRO:HA	1.53	0.89
1:C:95:ASP:HB3	1:C:109:PRO:HA	1.55	0.88
1:E:22:GLU:HB3	1:E:84:ILE:HG12	1.57	0.87
1:C:22:GLU:HB3	1:C:84:ILE:HG12	1.57	0.86
1:H:95:ASP:HB3	1:H:109:PRO:HA	1.55	0.86
1:H:22:GLU:HB3	1:H:84:ILE:HG12	1.56	0.86
1:D:22:GLU:HB3	1:D:84:ILE:HG12	1.58	0.85
1:A:22:GLU:HB3	1:A:84:ILE:HG12	1.58	0.83
1:B:95:ASP:HB3	1:B:109:PRO:HA	1.65	0.78
1:F:22:GLU:HB3	1:F:84:ILE:HG13	1.67	0.77
1:G:95:ASP:HB3	1:G:109:PRO:HA	1.65	0.77
1:E:9:LEU:HD12	1:E:9:LEU:O	1.84	0.76

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:28:LYS:HG3	1:B:50:ASN:HB3	1.68	0.75
1:F:95:ASP:HB3	1:F:109:PRO:HA	1.67	0.75
1:G:22:GLU:HB3	1:G:84:ILE:HG13	1.67	0.74
1:C:9:LEU:O	1:C:9:LEU:HD12	1.87	0.74
1:H:9:LEU:O	1:H:9:LEU:HD12	1.88	0.74
1:D:9:LEU:HD12	1:D:9:LEU:O	1.88	0.73
1:B:22:GLU:HB3	1:B:84:ILE:HG13	1.68	0.73
1:F:42:CYS:HB3	1:F:114:MET:HE3	1.70	0.73
1:G:42:CYS:HB3	1:G:114:MET:HE3	1.70	0.73
1:A:9:LEU:O	1:A:9:LEU:HD12	1.89	0.73
1:B:42:CYS:HB3	1:B:114:MET:HE3	1.71	0.73
1:B:57:THR:HG21	1:B:73:ARG:HH21	1.52	0.73
1:A:93:GLN:H	1:A:93:GLN:NE2	1.89	0.70
1:F:59:VAL:HG22	1:F:73:ARG:HG2	1.73	0.69
1:D:93:GLN:H	1:D:93:GLN:NE2	1.91	0.68
1:G:49:PHE:HA	1:G:56:ASN:OD1	1.94	0.67
1:B:49:PHE:HA	1:B:56:ASN:OD1	1.95	0.67
1:C:93:GLN:H	1:C:93:GLN:NE2	1.92	0.67
1:B:48:ARG:O	1:B:56:ASN:HA	1.95	0.66
1:E:93:GLN:NE2	1:E:93:GLN:H	1.94	0.66
1:F:48:ARG:O	1:F:56:ASN:HA	1.96	0.66
1:C:54:ASP:HB3	1:C:57:THR:CG2	2.26	0.65
1:H:93:GLN:H	1:H:93:GLN:NE2	1.94	0.65
1:E:54:ASP:HB3	1:E:57:THR:CG2	2.28	0.64
1:H:54:ASP:HB3	1:H:57:THR:CG2	2.28	0.64
1:C:7:SER:O	1:C:8:ASN:HB2	1.98	0.63
1:H:54:ASP:CG	1:H:73:ARG:HH12	2.06	0.63
1:A:7:SER:O	1:A:8:ASN:HB2	1.98	0.63
1:A:54:ASP:HB3	1:A:57:THR:CG2	2.29	0.63
1:G:49:PHE:O	1:G:50:ASN:HB2	1.97	0.63
1:G:83:SER:C	1:G:84:ILE:HD12	2.23	0.63
1:C:54:ASP:CG	1:C:73:ARG:HH12	2.07	0.62
1:G:48:ARG:O	1:G:56:ASN:HA	2.00	0.62
1:H:7:SER:O	1:H:8:ASN:HB2	1.99	0.62
1:A:44:HIS:NE2	2:I:2:GAL:H4	2.14	0.62
1:C:54:ASP:HB3	1:C:57:THR:HG21	1.81	0.62
1:H:54:ASP:HB3	1:H:57:THR:HG21	1.81	0.61
1:D:54:ASP:HB3	1:D:57:THR:CG2	2.31	0.61
1:A:54:ASP:HB3	1:A:57:THR:HG21	1.83	0.60
1:B:63:LYS:HD3	1:B:66:GLY:HA2	1.83	0.60
1:B:83:SER:C	1:B:84:ILE:HD12	2.26	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:83:SER:C	1:F:84:ILE:HD12	2.27	0.60
1:E:54:ASP:CG	1:E:73:ARG:HH12	2.08	0.60
1:C:66:GLY:O	1:G:67:THR:CG2	2.49	0.60
1:E:7:SER:O	1:E:8:ASN:HB2	2.02	0.59
1:D:54:ASP:HB3	1:D:57:THR:HG21	1.83	0.59
1:A:54:ASP:CG	1:A:73:ARG:HH12	2.11	0.58
1:D:7:SER:O	1:D:8:ASN:HB2	2.01	0.58
1:D:54:ASP:CG	1:D:73:ARG:HH12	2.10	0.58
1:E:76:ALA:HB1	1:E:104:HIS:CE1	2.39	0.58
1:B:10:ASN:HD21	1:B:12:LYS:HE2	1.69	0.58
1:G:63:LYS:HD3	1:G:66:GLY:HA2	1.85	0.57
1:C:66:GLY:O	1:G:67:THR:HG22	2.03	0.57
1:E:54:ASP:HB3	1:E:57:THR:HG21	1.85	0.57
1:F:63:LYS:HD3	1:F:66:GLY:HA2	1.85	0.57
1:C:76:ALA:HB1	1:C:104:HIS:CE1	2.40	0.57
1:A:76:ALA:HB1	1:A:104:HIS:CE1	2.40	0.56
1:H:76:ALA:HB1	1:H:104:HIS:CE1	2.40	0.56
1:E:30:PHE:HA	1:E:124:GLY:HA3	1.87	0.56
1:H:31:VAL:HB	1:H:123:ASP:HB3	1.88	0.55
1:F:43:LEU:HB2	1:F:96:LEU:HD22	1.87	0.55
1:B:59:VAL:HG13	1:B:71:GLU:HG3	1.89	0.55
1:B:43:LEU:HB2	1:B:96:LEU:HD22	1.88	0.55
1:C:26:ASP:O	1:C:28:LYS:HD2	2.07	0.55
1:D:93:GLN:H	1:D:93:GLN:CD	2.15	0.55
1:D:30:PHE:HA	1:D:124:GLY:HA3	1.88	0.54
1:E:26:ASP:O	1:E:28:LYS:HD2	2.07	0.54
1:F:49:PHE:HA	1:F:56:ASN:OD1	2.07	0.54
1:D:26:ASP:O	1:D:28:LYS:HD2	2.08	0.54
1:A:26:ASP:O	1:A:28:LYS:HD2	2.08	0.54
1:A:93:GLN:H	1:A:93:GLN:CD	2.15	0.54
1:C:54:ASP:OD1	1:C:73:ARG:NH1	2.41	0.54
1:F:48:ARG:C	1:F:50:ASN:H	2.13	0.54
1:F:10:ASN:HD21	1:F:12:LYS:HE2	1.73	0.54
1:G:43:LEU:HB2	1:G:96:LEU:HD22	1.90	0.54
1:C:9:LEU:O	1:C:118:ASN:HA	2.08	0.53
1:H:30:PHE:HA	1:H:124:GLY:HA3	1.89	0.53
1:A:9:LEU:O	1:A:118:ASN:HA	2.08	0.53
1:H:54:ASP:OD1	1:H:73:ARG:NH1	2.41	0.53
1:D:31:VAL:HB	1:D:123:ASP:HB3	1.90	0.53
1:D:76:ALA:HB1	1:D:104:HIS:CE1	2.42	0.53
1:E:88:CSO:HB3	1:E:99:LYS:HB3	1.90	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:31:VAL:HB	1:E:123:ASP:HB3	1.90	0.53
1:E:76:ALA:HB1	1:E:104:HIS:HE1	1.74	0.53
1:A:31:VAL:HB	1:A:123:ASP:HB3	1.90	0.53
1:E:47:PRO:HA	1:E:58:ILE:HA	1.91	0.53
1:F:18:LYS:HD3	1:F:86:GLU:OE1	2.08	0.53
1:H:93:GLN:H	1:H:93:GLN:CD	2.16	0.53
1:H:96:LEU:HD12	1:H:96:LEU:N	2.24	0.53
1:G:10:ASN:HD21	1:G:12:LYS:HE2	1.73	0.53
1:C:30:PHE:HA	1:C:124:GLY:HA3	1.91	0.53
1:C:31:VAL:HB	1:C:123:ASP:HB3	1.90	0.53
1:C:93:GLN:H	1:C:93:GLN:CD	2.17	0.52
1:D:47:PRO:HA	1:D:58:ILE:HA	1.91	0.52
1:D:88:CSO:HB3	1:D:99:LYS:HB3	1.91	0.52
1:E:93:GLN:H	1:E:93:GLN:CD	2.17	0.52
1:A:47:PRO:HA	1:A:58:ILE:HA	1.91	0.52
1:D:96:LEU:N	1:D:96:LEU:HD12	2.25	0.52
1:H:26:ASP:O	1:H:28:LYS:HD2	2.09	0.52
1:B:18:LYS:HD3	1:B:86:GLU:OE1	2.09	0.52
1:H:47:PRO:HA	1:H:58:ILE:HA	1.90	0.52
1:C:47:PRO:HA	1:C:58:ILE:HA	1.91	0.52
1:E:96:LEU:N	1:E:96:LEU:HD12	2.25	0.52
1:C:31:VAL:HG12	1:C:32:LEU:N	2.25	0.52
1:E:98:ILE:N	1:E:98:ILE:HD12	2.25	0.52
1:G:90:THR:HG22	1:G:91:PHE:N	2.24	0.52
1:H:98:ILE:HD12	1:H:98:ILE:N	2.25	0.52
1:A:66:GLY:O	1:E:67:THR:HB	2.10	0.52
1:B:73:ARG:C	1:B:74:GLU:HG2	2.35	0.52
1:H:9:LEU:O	1:H:118:ASN:HA	2.09	0.52
1:A:90:THR:HG22	1:A:91:PHE:N	2.25	0.51
1:A:96:LEU:HD12	1:A:96:LEU:N	2.25	0.51
1:D:54:ASP:OD1	1:D:73:ARG:NH1	2.43	0.51
1:E:91:PHE:CE2	1:E:117:ILE:HD13	2.46	0.51
1:H:90:THR:HG22	1:H:91:PHE:N	2.25	0.51
1:C:96:LEU:HD12	1:C:96:LEU:N	2.26	0.51
1:G:92:ASP:C	1:G:92:ASP:OD2	2.54	0.51
1:F:100:LEU:HD12	1:F:104:HIS:HD2	1.76	0.51
1:A:30:PHE:HA	1:A:124:GLY:HA3	1.92	0.50
1:C:76:ALA:HB1	1:C:104:HIS:HE1	1.76	0.50
1:G:18:LYS:HD3	1:G:86:GLU:OE1	2.10	0.50
1:B:92:ASP:OD2	1:B:92:ASP:C	2.54	0.50
1:C:10:ASN:HA	1:C:118:ASN:ND2	2.27	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:90:THR:HG22	1:F:91:PHE:N	2.26	0.50
1:H:42:CYS:HB3	1:H:114:MET:HE3	1.94	0.50
1:C:42:CYS:HB3	1:C:114:MET:HE3	1.93	0.50
1:D:90:THR:HG22	1:D:91:PHE:N	2.26	0.50
1:E:90:THR:HG22	1:E:91:PHE:N	2.27	0.50
1:C:98:ILE:HD12	1:C:98:ILE:N	2.27	0.49
1:F:4:LEU:C	1:F:4:LEU:HD23	2.37	0.49
1:F:48:ARG:HB3	1:F:50:ASN:C	2.36	0.49
1:H:76:ALA:HB1	1:H:104:HIS:HE1	1.75	0.49
1:A:54:ASP:OD1	1:A:73:ARG:NH1	2.42	0.49
1:E:42:CYS:HB3	1:E:114:MET:HE3	1.93	0.49
1:F:92:ASP:OD2	1:F:92:ASP:C	2.55	0.49
1:A:43:LEU:HB2	1:A:96:LEU:HD22	1.94	0.49
1:C:90:THR:HG22	1:C:91:PHE:N	2.26	0.49
1:D:42:CYS:HB3	1:D:114:MET:HE3	1.94	0.49
1:D:9:LEU:O	1:D:118:ASN:HA	2.12	0.49
1:F:30:PHE:CZ	1:F:47:PRO:HG3	2.47	0.49
1:D:91:PHE:CE2	1:D:117:ILE:HD13	2.48	0.49
1:A:76:ALA:HB1	1:A:104:HIS:HE1	1.77	0.49
1:A:100:LEU:HB3	1:A:101:PRO:HD2	1.95	0.49
1:B:49:PHE:N	1:B:49:PHE:CD2	2.78	0.49
1:B:100:LEU:HD12	1:B:104:HIS:HD2	1.78	0.49
1:E:50:ASN:CG	1:H:25:PRO:HG3	2.38	0.49
1:H:10:ASN:HA	1:H:118:ASN:ND2	2.28	0.49
1:H:88:CSO:HB3	1:H:99:LYS:HB3	1.95	0.49
1:B:4:LEU:C	1:B:4:LEU:HD23	2.38	0.49
1:A:10:ASN:HA	1:A:118:ASN:ND2	2.27	0.48
1:A:93:GLN:HG3	1:A:115:GLU:OE2	2.13	0.48
1:C:43:LEU:HB2	1:C:96:LEU:HD22	1.95	0.48
1:D:43:LEU:HB2	1:D:96:LEU:HD22	1.95	0.48
1:E:50:ASN:OD1	1:H:25:PRO:HG3	2.12	0.48
1:E:54:ASP:OD1	1:E:73:ARG:NH1	2.44	0.48
1:C:9:LEU:HB3	1:D:4:LEU:HD13	1.95	0.48
1:E:43:LEU:HB2	1:E:96:LEU:HD22	1.95	0.48
1:A:31:VAL:HG12	1:A:32:LEU:N	2.29	0.48
1:F:64:ASP:O	1:F:65:ASP:C	2.56	0.48
1:F:88:CSO:HB3	1:F:99:LYS:HB3	1.95	0.48
1:D:93:GLN:NE2	1:D:93:GLN:N	2.61	0.48
1:E:9:LEU:O	1:E:118:ASN:HA	2.14	0.48
1:F:23:LEU:N	1:F:23:LEU:HD22	2.29	0.48
1:H:43:LEU:HB2	1:H:96:LEU:HD22	1.96	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:64:ASP:O	1:B:65:ASP:C	2.55	0.48
1:D:76:ALA:HB1	1:D:104:HIS:HE1	1.78	0.48
1:A:88:CSO:HB3	1:A:99:LYS:HB3	1.96	0.48
1:A:91:PHE:CE2	1:A:117:ILE:HD13	2.49	0.48
1:B:90:THR:HG22	1:B:91:PHE:N	2.28	0.48
1:A:39:ASN:OD1	1:E:66:GLY:HA3	2.14	0.47
1:D:98:ILE:HD12	1:D:98:ILE:N	2.28	0.47
1:H:91:PHE:CE2	1:H:117:ILE:HD13	2.49	0.47
1:C:93:GLN:HG3	1:C:115:GLU:OE2	2.15	0.47
1:G:30:PHE:CZ	1:G:47:PRO:HG3	2.49	0.47
1:B:88:CSO:HB3	1:B:99:LYS:HB3	1.97	0.47
1:G:23:LEU:HD22	1:G:23:LEU:N	2.29	0.47
1:E:31:VAL:HG12	1:E:32:LEU:N	2.29	0.47
1:G:100:LEU:HD12	1:G:104:HIS:HD2	1.78	0.47
1:A:4:LEU:CD1	1:B:9:LEU:HB3	2.45	0.47
1:A:93:GLN:HG3	1:A:115:GLU:CD	2.40	0.47
1:G:64:ASP:O	1:G:65:ASP:C	2.58	0.47
1:A:42:CYS:HB3	1:A:114:MET:HE3	1.96	0.47
1:H:93:GLN:HG3	1:H:115:GLU:OE2	2.15	0.46
1:A:98:ILE:HD12	1:A:98:ILE:N	2.29	0.46
1:C:88:CSO:HB3	1:C:99:LYS:HB3	1.97	0.46
1:C:91:PHE:CE2	1:C:117:ILE:HD13	2.50	0.46
1:E:10:ASN:HA	1:E:118:ASN:ND2	2.29	0.46
1:A:4:LEU:HD13	1:B:9:LEU:HB3	1.96	0.46
1:B:10:ASN:ND2	1:B:12:LYS:HE2	2.30	0.46
1:A:4:LEU:HD12	1:B:7:SER:O	2.15	0.46
1:D:93:GLN:HG3	1:D:115:GLU:OE2	2.16	0.46
1:B:30:PHE:CZ	1:B:47:PRO:HG3	2.50	0.46
1:D:47:PRO:HB3	1:D:58:ILE:HG12	1.98	0.46
1:H:31:VAL:HG12	1:H:32:LEU:N	2.30	0.46
1:D:10:ASN:HA	1:D:118:ASN:ND2	2.31	0.46
1:C:93:GLN:NE2	1:C:93:GLN:N	2.63	0.46
1:G:96:LEU:N	1:G:96:LEU:HD12	2.31	0.46
1:A:47:PRO:HB3	1:A:58:ILE:HG12	1.98	0.46
1:D:31:VAL:HG12	1:D:32:LEU:N	2.30	0.46
1:B:23:LEU:HD22	1:B:23:LEU:N	2.30	0.46
1:H:54:ASP:CG	1:H:73:ARG:NH1	2.73	0.45
1:H:93:GLN:HG3	1:H:115:GLU:CD	2.41	0.45
1:E:93:GLN:HG3	1:E:115:GLU:OE2	2.16	0.45
1:G:4:LEU:C	1:G:4:LEU:HD23	2.41	0.45
1:G:102:ASP:OD1	1:G:104:HIS:HB3	2.17	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:93:GLN:HG3	1:D:115:GLU:CD	2.41	0.45
1:G:100:LEU:HB3	1:G:101:PRO:HD2	1.99	0.45
1:G:59:VAL:HG22	1:G:73:ARG:HG2	1.97	0.45
1:F:30:PHE:HB2	1:F:126:PHE:HB2	1.99	0.45
1:H:11:LEU:HD21	1:H:133:PHE:HD2	1.82	0.45
1:F:102:ASP:OD1	1:F:104:HIS:HB3	2.17	0.45
1:F:81:PRO:C	1:F:83:SER:H	2.24	0.44
1:G:88:CSO:HB3	1:G:99:LYS:HB3	1.99	0.44
1:A:90:THR:HG22	1:A:91:PHE:H	1.81	0.44
1:C:54:ASP:CG	1:C:73:ARG:NH1	2.74	0.44
1:G:30:PHE:HB2	1:G:126:PHE:HB2	1.99	0.44
1:A:93:GLN:NE2	1:A:93:GLN:N	2.61	0.44
1:B:96:LEU:N	1:B:96:LEU:HD12	2.33	0.44
1:D:91:PHE:CD2	1:D:117:ILE:HD13	2.53	0.44
1:E:54:ASP:CG	1:E:73:ARG:NH1	2.75	0.44
1:F:16:CYS:HB3	1:F:134:GLU:HB2	2.00	0.44
1:H:47:PRO:HB3	1:H:58:ILE:HG12	1.99	0.44
1:A:11:LEU:HB2	1:A:117:ILE:HB	1.99	0.44
1:C:11:LEU:HD21	1:C:133:PHE:HD2	1.82	0.44
1:E:11:LEU:HB2	1:E:117:ILE:HB	1.99	0.44
1:B:81:PRO:C	1:B:83:SER:H	2.26	0.43
1:E:47:PRO:HB3	1:E:58:ILE:HG12	2.00	0.43
1:E:93:GLN:HG3	1:E:115:GLU:CD	2.43	0.43
1:B:30:PHE:HB2	1:B:126:PHE:HB2	2.00	0.43
1:E:4:LEU:CD1	1:F:9:LEU:HB3	2.48	0.43
1:F:96:LEU:HD12	1:F:96:LEU:N	2.34	0.43
1:D:67:THR:HG22	1:H:66:GLY:O	2.18	0.43
1:D:11:LEU:HD21	1:D:133:PHE:HD2	1.83	0.43
1:D:33:ASN:O	1:D:120:MET:HA	2.19	0.43
1:E:50:ASN:HB2	1:H:25:PRO:HG2	2.01	0.43
1:F:28:LYS:HG3	1:F:50:ASN:CB	2.48	0.43
1:G:81:PRO:C	1:G:83:SER:H	2.26	0.43
1:H:11:LEU:HB2	1:H:117:ILE:HB	1.99	0.43
1:F:9:LEU:HD12	1:F:9:LEU:O	2.18	0.43
1:A:122:ALA:HB2	1:A:128:ILE:HD11	2.01	0.43
1:C:11:LEU:HB2	1:C:117:ILE:HB	2.01	0.43
1:C:100:LEU:HB3	1:C:101:PRO:HD2	2.00	0.43
1:E:50:ASN:HB2	1:H:25:PRO:CG	2.49	0.43
1:A:10:ASN:O	1:A:10:ASN:CG	2.62	0.43
1:D:32:LEU:HD12	1:D:32:LEU:HA	1.85	0.43
1:C:47:PRO:HB3	1:C:58:ILE:HG12	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:93:GLN:HG3	1:C:115:GLU:CD	2.43	0.42
1:F:100:LEU:HB3	1:F:101:PRO:HD2	2.00	0.42
1:H:33:ASN:O	1:H:120:MET:HA	2.19	0.42
1:H:93:GLN:NE2	1:H:93:GLN:N	2.64	0.42
1:C:90:THR:HG22	1:C:91:PHE:H	1.83	0.42
1:C:31:VAL:CG1	1:C:32:LEU:N	2.82	0.42
1:D:90:THR:HG22	1:D:91:PHE:H	1.83	0.42
1:E:91:PHE:CD2	1:E:117:ILE:HD13	2.54	0.42
1:A:32:LEU:HD12	1:A:32:LEU:HA	1.92	0.42
1:A:44:HIS:NE2	2:I:2:GAL:C4	2.81	0.42
1:B:71:GLU:O	1:B:71:GLU:HG2	2.19	0.42
1:C:33:ASN:O	1:C:120:MET:HA	2.19	0.42
1:D:11:LEU:HB2	1:D:117:ILE:HB	2.00	0.42
1:E:33:ASN:O	1:E:120:MET:HA	2.19	0.42
1:F:10:ASN:ND2	1:F:12:LYS:HE2	2.35	0.42
1:B:100:LEU:HB3	1:B:101:PRO:HD2	2.01	0.42
1:E:10:ASN:CG	1:E:10:ASN:O	2.63	0.42
1:F:84:ILE:O	1:F:84:ILE:HG22	2.20	0.42
1:H:90:THR:HG22	1:H:91:PHE:H	1.85	0.42
1:D:54:ASP:CG	1:D:73:ARG:NH1	2.77	0.42
1:G:9:LEU:HD12	1:G:9:LEU:O	2.20	0.42
1:E:26:ASP:OD1	1:E:26:ASP:N	2.53	0.41
1:G:59:VAL:HG13	1:G:71:GLU:HG3	2.02	0.41
1:E:90:THR:HG22	1:E:91:PHE:H	1.84	0.41
1:E:93:GLN:NE2	1:E:93:GLN:N	2.64	0.41
1:F:22:GLU:HA	1:F:84:ILE:HA	2.00	0.41
1:A:11:LEU:HD21	1:A:133:PHE:HD2	1.85	0.41
1:E:11:LEU:HD21	1:E:133:PHE:HD2	1.85	0.41
1:H:100:LEU:HB3	1:H:101:PRO:HD2	2.01	0.41
1:C:10:ASN:O	1:C:10:ASN:CG	2.64	0.41
1:B:22:GLU:HA	1:B:84:ILE:HA	2.02	0.41
1:A:3:GLY:HA3	1:A:122:ALA:O	2.21	0.41
1:B:67:THR:HG22	1:F:66:GLY:O	2.20	0.41
1:D:122:ALA:HB2	1:D:128:ILE:HD11	2.03	0.41
1:F:81:PRO:C	1:F:83:SER:N	2.79	0.41
1:H:26:ASP:OD1	1:H:26:ASP:N	2.54	0.41
1:H:91:PHE:CD2	1:H:117:ILE:HD13	2.56	0.41
1:H:122:ALA:HB2	1:H:128:ILE:HD11	2.03	0.41
1:B:9:LEU:HD12	1:B:9:LEU:O	2.21	0.40
1:C:122:ALA:HB2	1:C:128:ILE:HD11	2.04	0.40
1:H:10:ASN:O	1:H:10:ASN:CG	2.63	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:31:VAL:CG1	1:A:32:LEU:N	2.85	0.40
1:C:66:GLY:CA	1:G:67:THR:HG22	2.51	0.40
1:E:100:LEU:HB3	1:E:101:PRO:HD2	2.03	0.40
1:F:23:LEU:HD21	1:F:85:THR:OG1	2.20	0.40
1:A:26:ASP:OD1	1:A:26:ASP:N	2.54	0.40
1:C:23:LEU:HD11	1:C:30:PHE:CE2	2.56	0.40
1:E:32:LEU:HD12	1:E:32:LEU:HA	1.89	0.40
1:G:30:PHE:CB	1:G:126:PHE:HB2	2.52	0.40
1:H:23:LEU:HD11	1:H:30:PHE:CE2	2.57	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	128/134 (96%)	119 (93%)	8 (6%)	1 (1%)	16	37
1	B	114/134 (85%)	109 (96%)	2 (2%)	3 (3%)	4	12
1	C	128/134 (96%)	120 (94%)	7 (6%)	1 (1%)	16	37
1	D	128/134 (96%)	119 (93%)	8 (6%)	1 (1%)	16	37
1	E	128/134 (96%)	120 (94%)	7 (6%)	1 (1%)	16	37
1	F	114/134 (85%)	108 (95%)	5 (4%)	1 (1%)	14	34
1	G	114/134 (85%)	107 (94%)	6 (5%)	1 (1%)	14	34
1	H	128/134 (96%)	120 (94%)	7 (6%)	1 (1%)	16	37
All	All	982/1072 (92%)	922 (94%)	50 (5%)	10 (1%)	12	32

All (10) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	10	ASN

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Mol	Chain	Res	Type
1	A	94	ALA
1	B	82	GLY
1	C	94	ALA
1	D	94	ALA
1	E	94	ALA
1	F	10	ASN
1	G	10	ASN
1	H	94	ALA
1	B	65	ASP

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	109/110 (99%)	102 (94%)	7 (6%)	16	38
1	B	102/110 (93%)	96 (94%)	6 (6%)	18	41
1	C	109/110 (99%)	102 (94%)	7 (6%)	16	38
1	D	109/110 (99%)	102 (94%)	7 (6%)	16	38
1	E	109/110 (99%)	102 (94%)	7 (6%)	16	38
1	F	102/110 (93%)	96 (94%)	6 (6%)	18	41
1	G	102/110 (93%)	96 (94%)	6 (6%)	18	41
1	H	109/110 (99%)	102 (94%)	7 (6%)	16	38
All	All	851/880 (97%)	798 (94%)	53 (6%)	16	39

All (53) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	18	LYS
1	A	57	THR
1	A	67	THR
1	A	93	GLN
1	A	95	ASP
1	A	113	ASN

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Mol	Chain	Res	Type
1	A	115	GLU
1	B	9	LEU
1	B	11	LEU
1	B	49	PHE
1	B	70	THR
1	B	95	ASP
1	B	113	ASN
1	C	18	LYS
1	C	57	THR
1	C	67	THR
1	C	93	GLN
1	C	95	ASP
1	C	113	ASN
1	C	115	GLU
1	D	18	LYS
1	D	57	THR
1	D	67	THR
1	D	93	GLN
1	D	95	ASP
1	D	113	ASN
1	D	115	GLU
1	E	18	LYS
1	E	57	THR
1	E	67	THR
1	E	93	GLN
1	E	95	ASP
1	E	113	ASN
1	E	115	GLU
1	F	9	LEU
1	F	11	LEU
1	F	49	PHE
1	F	95	ASP
1	F	113	ASN
1	F	134	GLU
1	G	9	LEU
1	G	11	LEU
1	G	74	GLU
1	G	95	ASP
1	G	113	ASN
1	G	134	GLU
1	H	18	LYS
1	H	57	THR

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Mol	Chain	Res	Type
1	H	67	THR
1	H	93	GLN
1	H	95	ASP
1	H	113	ASN
1	H	115	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (46) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	50	ASN
1	A	52	HIS
1	A	56	ASN
1	A	93	GLN
1	A	113	ASN
1	A	118	ASN
1	B	10	ASN
1	B	39	ASN
1	B	72	GLN
1	B	104	HIS
1	B	113	ASN
1	C	50	ASN
1	C	56	ASN
1	C	93	GLN
1	C	104	HIS
1	C	113	ASN
1	C	118	ASN
1	D	50	ASN
1	D	52	HIS
1	D	56	ASN
1	D	93	GLN
1	D	113	ASN
1	D	118	ASN
1	E	8	ASN
1	E	39	ASN
1	E	50	ASN
1	E	56	ASN
1	E	93	GLN
1	E	104	HIS
1	E	113	ASN
1	E	118	ASN
1	F	10	ASN
1	F	104	HIS

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Mol	Chain	Res	Type
1	F	113	ASN
1	G	10	ASN
1	G	39	ASN
1	G	50	ASN
1	G	72	GLN
1	G	104	HIS
1	G	113	ASN
1	H	50	ASN
1	H	56	ASN
1	H	93	GLN
1	H	104	HIS
1	H	113	ASN
1	H	118	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

16 non-standard protein/DNA/RNA residues 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	$\# Z > 2$
1	CSO	H	130	1	3,6,7	0.69	0	1,6,8	0.33	0
1	CSO	D	130	1	3,6,7	0.70	0	1,6,8	0.26	0
1	CSO	D	88	1	3,6,7	0.70	0	1,6,8	0.16	0
1	CSO	B	130	1	3,6,7	0.72	0	1,6,8	0.53	0
1	CSO	A	130	1	3,6,7	0.76	0	1,6,8	0.25	0
1	CSO	E	88	1	3,6,7	0.72	0	1,6,8	0.00	0
1	CSO	F	88	1	3,6,7	0.86	0	1,6,8	0.08	0
1	CSO	F	130	1	3,6,7	0.68	0	1,6,8	0.55	0
1	CSO	C	88	1	3,6,7	0.71	0	1,6,8	0.15	0
1	CSO	B	88	1	3,6,7	1.00	0	1,6,8	0.07	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
1	CSO	A	88	1	3,6,7	0.82	0	1,6,8	0.20	0
1	CSO	H	88	1	3,6,7	0.74	0	1,6,8	0.01	0
1	CSO	G	88	1	3,6,7	0.82	0	1,6,8	0.08	0
1	CSO	E	130	1	3,6,7	0.65	0	1,6,8	0.31	0
1	CSO	G	130	1	3,6,7	0.72	0	1,6,8	0.27	0
1	CSO	C	130	1	3,6,7	0.75	0	1,6,8	0.12	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
1	CSO	H	130	1	-	1/1/5/7	-
1	CSO	D	130	1	-	1/1/5/7	-
1	CSO	D	88	1	-	0/1/5/7	-
1	CSO	B	130	1	-	0/1/5/7	-
1	CSO	A	130	1	-	1/1/5/7	-
1	CSO	E	88	1	-	0/1/5/7	-
1	CSO	F	88	1	-	0/1/5/7	-
1	CSO	F	130	1	-	0/1/5/7	-
1	CSO	C	88	1	-	0/1/5/7	-
1	CSO	B	88	1	-	0/1/5/7	-
1	CSO	A	88	1	-	0/1/5/7	-
1	CSO	H	88	1	-	0/1/5/7	-
1	CSO	G	88	1	-	0/1/5/7	-
1	CSO	E	130	1	-	1/1/5/7	-
1	CSO	G	130	1	-	0/1/5/7	-
1	CSO	C	130	1	-	1/1/5/7	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (5) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	A	130	CSO	N-CA-CB-SG
1	C	130	CSO	N-CA-CB-SG
1	D	130	CSO	N-CA-CB-SG
1	E	130	CSO	N-CA-CB-SG
1	H	130	CSO	N-CA-CB-SG

There are no ring outliers.

8 monomers are involved in 8 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	D	88	CSO	1	0
1	E	88	CSO	1	0
1	F	88	CSO	1	0
1	C	88	CSO	1	0
1	B	88	CSO	1	0
1	A	88	CSO	1	0
1	H	88	CSO	1	0
1	G	88	CSO	1	0

5.5 Carbohydrates

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	$\# Z > 2$
2	BGC	I	1	2	12,12,12	0.53	0	17,17,17	0.50	0
2	GAL	I	2	2	11,11,12	0.52	0	15,15,17	0.56	0
2	BGC	J	1	2	12,12,12	0.37	0	17,17,17	0.59	0
2	GAL	J	2	2	11,11,12	0.51	0	15,15,17	0.56	0
2	BGC	K	1	2	12,12,12	0.45	0	17,17,17	0.38	0
2	GAL	K	2	2	11,11,12	0.41	0	15,15,17	0.42	0
2	BGC	L	1	2	12,12,12	0.37	0	17,17,17	0.51	0
2	GAL	L	2	2	11,11,12	0.38	0	15,15,17	0.61	0
2	BGC	M	1	2	12,12,12	0.42	0	17,17,17	0.42	0
2	GAL	M	2	2	11,11,12	0.54	0	15,15,17	0.38	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
2	BGC	I	1	2	-	2/2/22/22	0/1/1/1
2	GAL	I	2	2	-	2/2/19/22	0/1/1/1
2	BGC	J	1	2	-	0/2/22/22	0/1/1/1
2	GAL	J	2	2	-	1/2/19/22	0/1/1/1
2	BGC	K	1	2	-	0/2/22/22	0/1/1/1
2	GAL	K	2	2	-	0/2/19/22	0/1/1/1
2	BGC	L	1	2	-	2/2/22/22	0/1/1/1
2	GAL	L	2	2	-	2/2/19/22	0/1/1/1
2	BGC	M	1	2	-	0/2/22/22	0/1/1/1
2	GAL	M	2	2	-	1/2/19/22	0/1/1/1

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (10) torsion outliers are listed below:

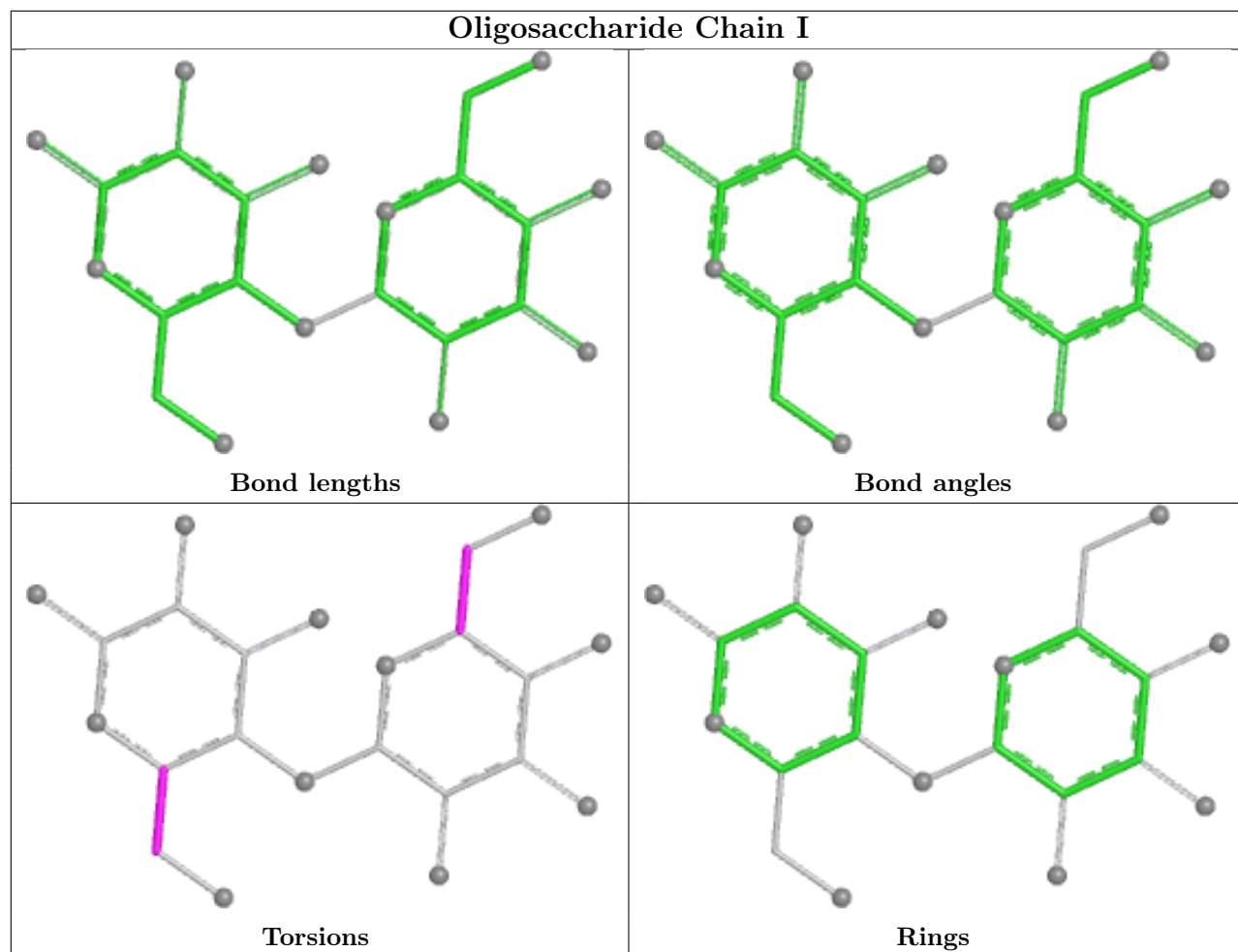
Mol	Chain	Res	Type	Atoms
2	I	1	BGC	O5-C5-C6-O6
2	I	1	BGC	C4-C5-C6-O6
2	I	2	GAL	O5-C5-C6-O6
2	L	1	BGC	C4-C5-C6-O6
2	I	2	GAL	C4-C5-C6-O6
2	L	1	BGC	O5-C5-C6-O6
2	L	2	GAL	O5-C5-C6-O6
2	M	2	GAL	C4-C5-C6-O6
2	J	2	GAL	C4-C5-C6-O6
2	L	2	GAL	C4-C5-C6-O6

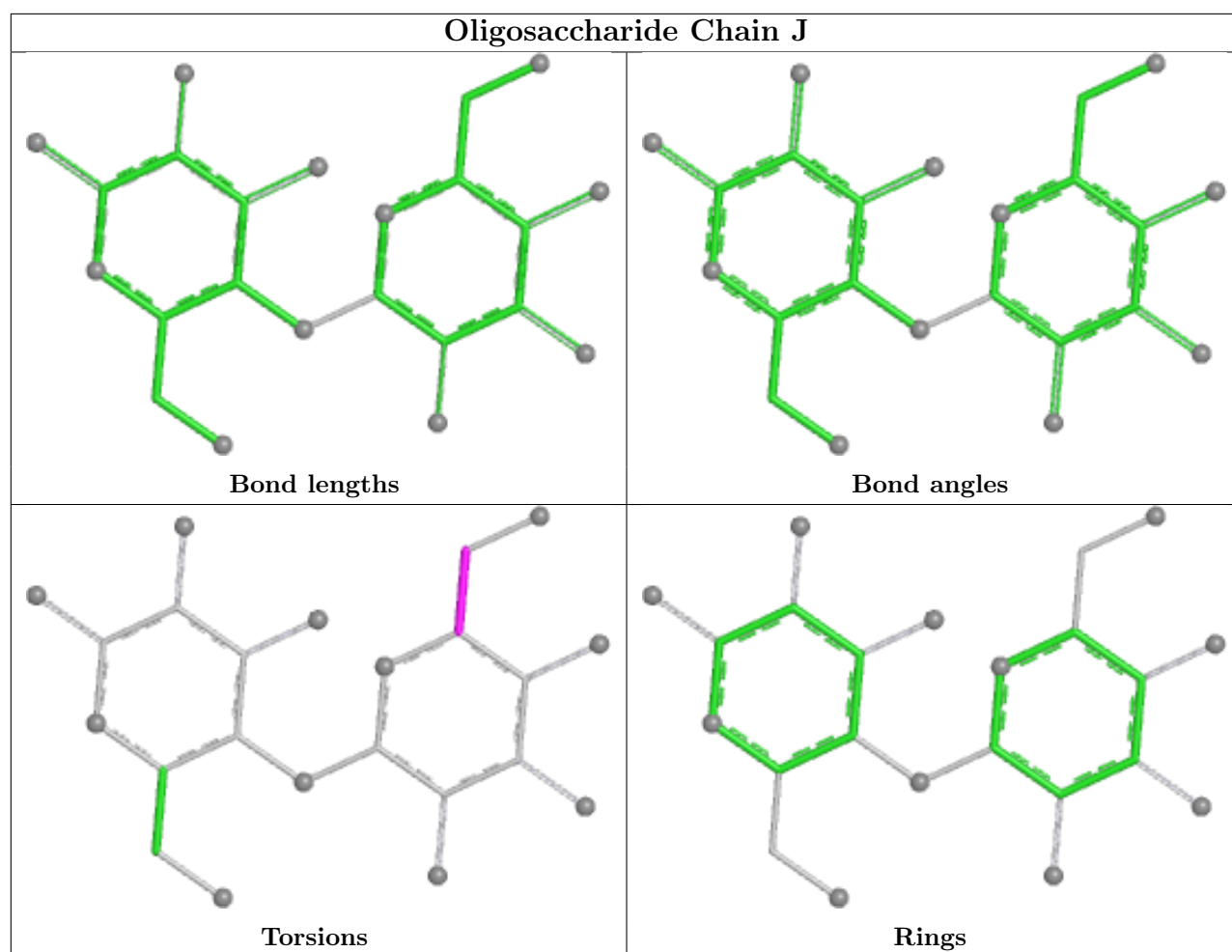
There are no ring outliers.

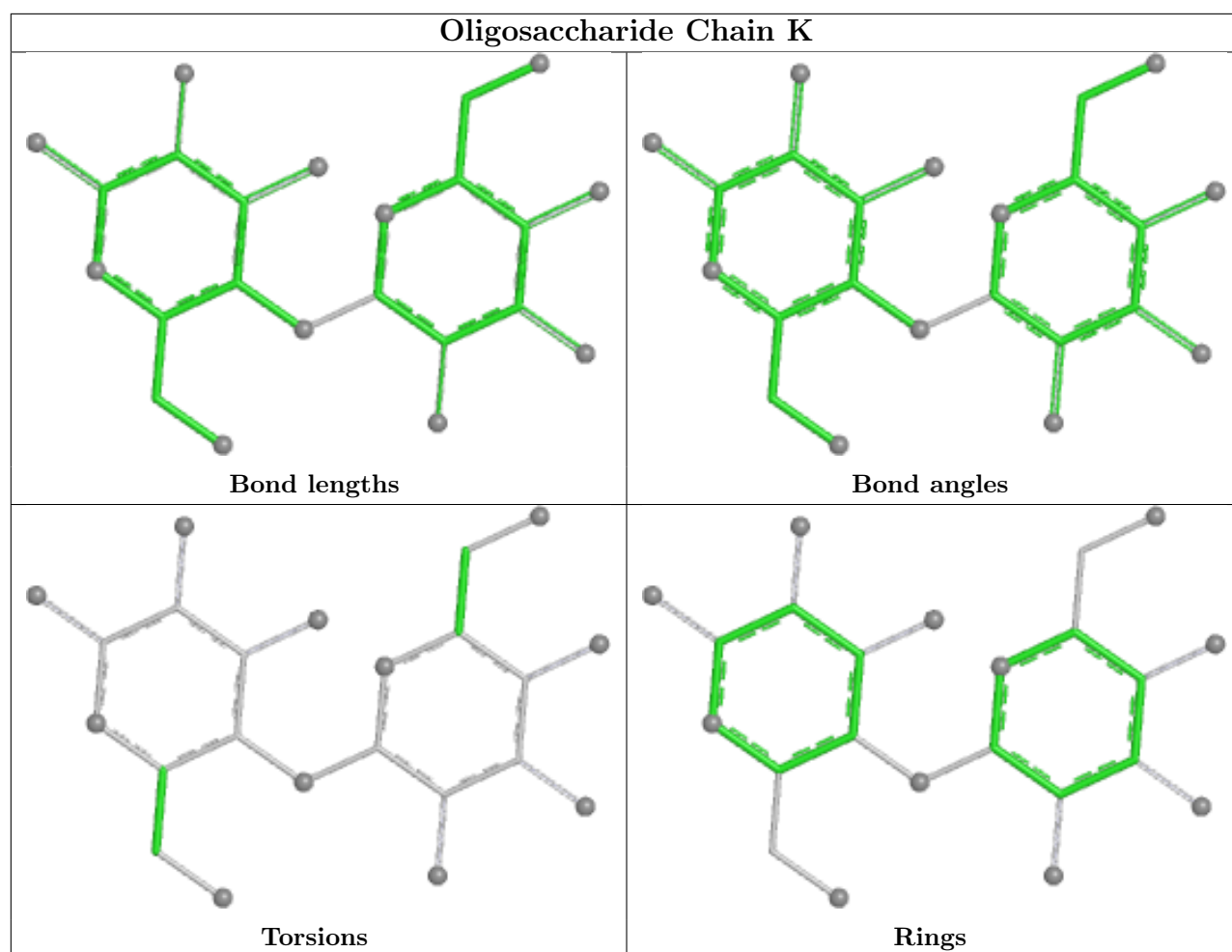
1 monomer is involved in 2 short contacts:

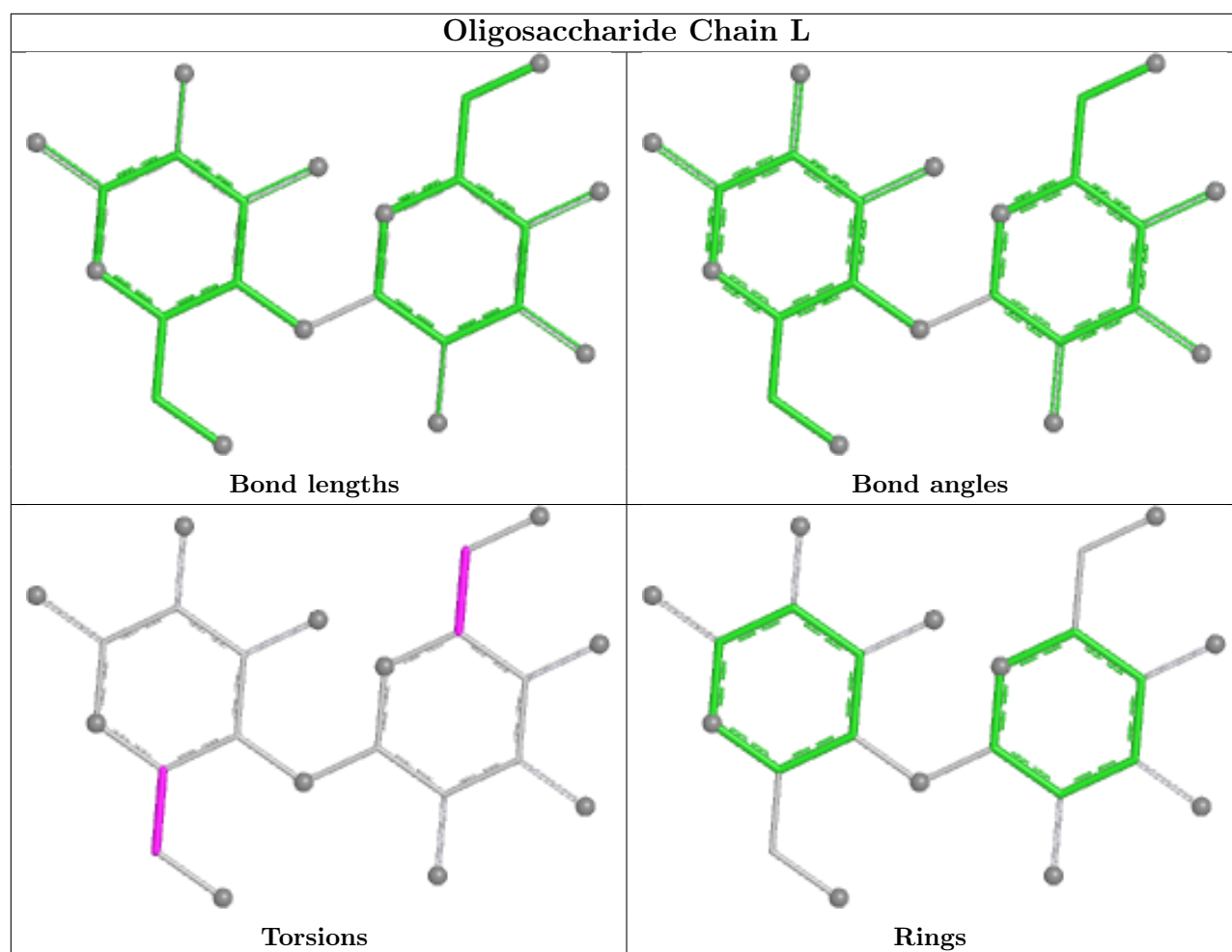
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	I	2	GAL	2	0

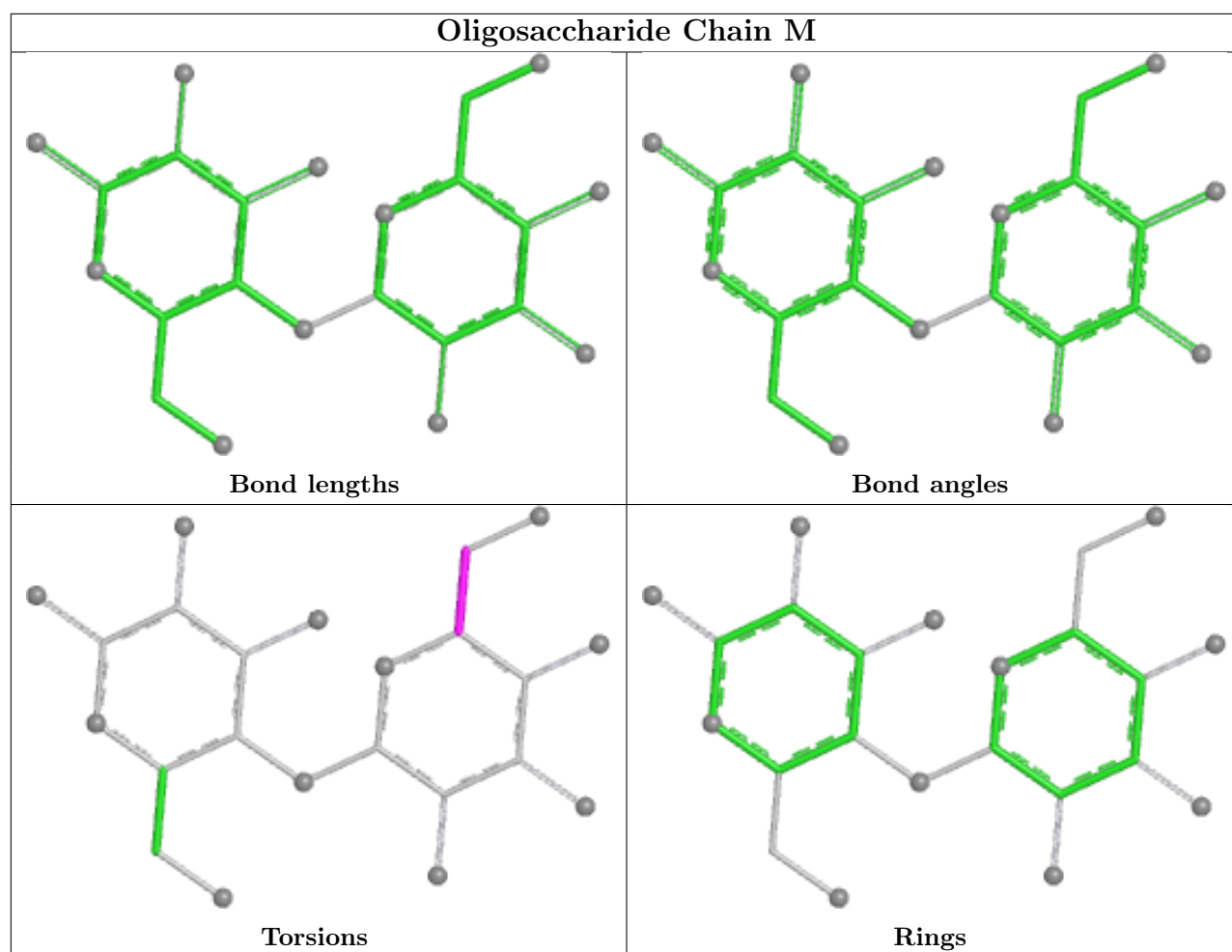
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for oligosaccharide.











5.6 Ligand geometry [i](#)

There are no ligands in this entry.

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	130/134 (97%)	0.05	1 (0%) 82 78	14, 28, 42, 50	0
1	B	120/134 (89%)	-0.12	0 100 100	15, 27, 43, 65	0
1	C	130/134 (97%)	-0.05	0 100 100	13, 28, 40, 49	0
1	D	130/134 (97%)	0.04	1 (0%) 82 78	14, 27, 40, 48	0
1	E	130/134 (97%)	0.01	2 (1%) 72 65	12, 27, 40, 48	0
1	F	120/134 (89%)	-0.05	2 (1%) 69 62	15, 27, 43, 63	0
1	G	120/134 (89%)	-0.19	0 100 100	14, 27, 41, 46	0
1	H	130/134 (97%)	-0.03	0 100 100	12, 27, 40, 48	0
All	All	1010/1072 (94%)	-0.04	6 (0%) 85 82	12, 27, 41, 65	0

All (6) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	F	50	ASN	4.6
1	D	112	LEU	2.9
1	E	115	GLU	2.3
1	E	93	GLN	2.2
1	F	115	GLU	2.1
1	A	54	ASP	2.0

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

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
1	CSO	A	130	7/8	0.87	0.10	20,23,28,30	0
1	CSO	G	130	7/8	0.92	0.08	14,18,21,23	0
1	CSO	C	130	7/8	0.93	0.09	20,21,26,26	0
1	CSO	C	88	7/8	0.93	0.08	20,21,22,23	0
1	CSO	D	130	7/8	0.94	0.08	20,20,23,23	0
1	CSO	F	88	7/8	0.94	0.09	11,18,19,19	0
1	CSO	B	130	7/8	0.94	0.06	18,20,23,24	0
1	CSO	E	88	7/8	0.95	0.07	18,21,22,23	0
1	CSO	E	130	7/8	0.95	0.07	19,20,23,25	0
1	CSO	D	88	7/8	0.95	0.08	18,20,21,21	0
1	CSO	F	130	7/8	0.95	0.08	18,20,21,23	0
1	CSO	A	88	7/8	0.95	0.07	21,21,22,25	0
1	CSO	H	88	7/8	0.95	0.08	20,20,20,21	0
1	CSO	H	130	7/8	0.95	0.06	19,20,23,25	0
1	CSO	G	88	7/8	0.96	0.08	13,16,17,17	0
1	CSO	B	88	7/8	0.97	0.07	11,16,16,17	0

6.3 Carbohydrates

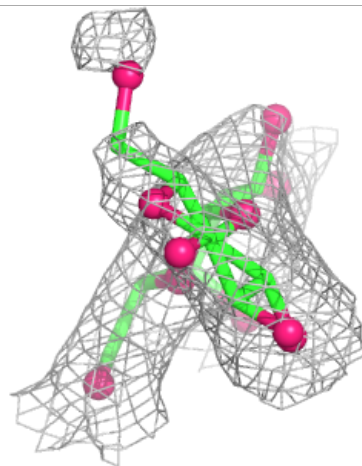
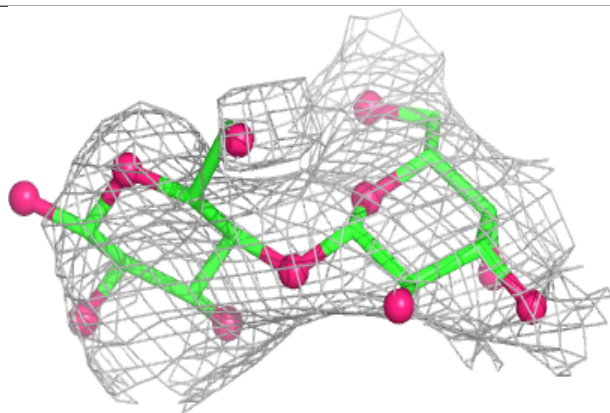
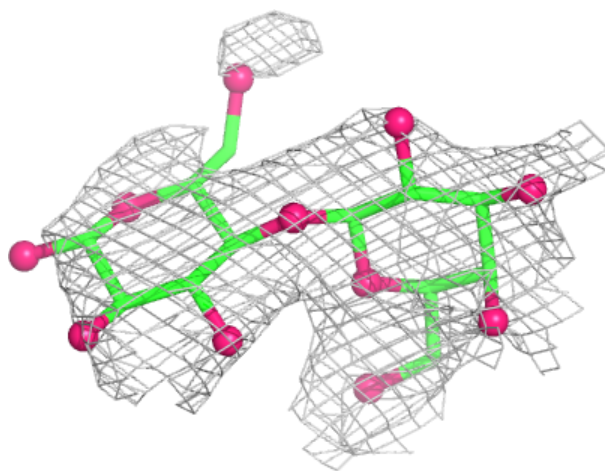
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	BGC	I	1	12/12	0.79	0.14	26,27,28,30	12
2	BGC	K	1	12/12	0.86	0.12	28,29,29,29	12
2	BGC	M	1	12/12	0.87	0.10	36,38,40,41	0
2	BGC	L	1	12/12	0.88	0.12	26,28,30,32	0
2	GAL	J	2	11/12	0.90	0.10	28,30,31,32	0
2	GAL	I	2	11/12	0.90	0.11	19,24,26,27	11
2	BGC	J	1	12/12	0.91	0.09	28,29,30,31	0
2	GAL	K	2	11/12	0.91	0.11	30,30,31,31	11
2	GAL	L	2	11/12	0.95	0.07	26,30,32,33	0
2	GAL	M	2	11/12	0.96	0.07	26,33,35,36	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.

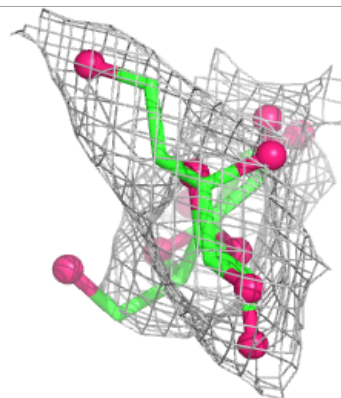
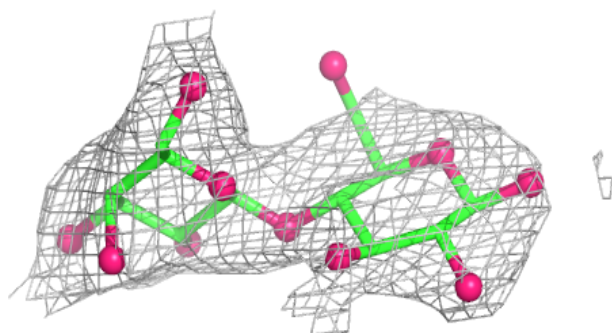
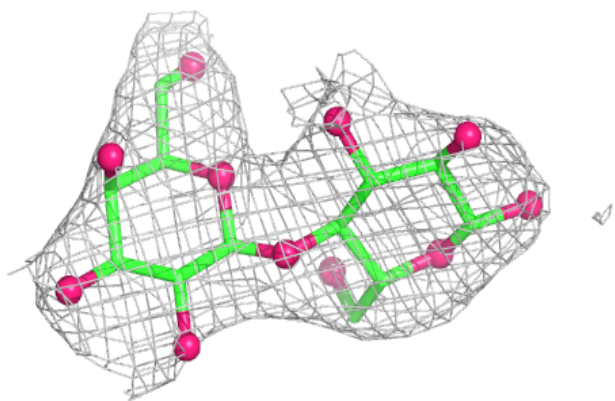
Electron density around Chain I:

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

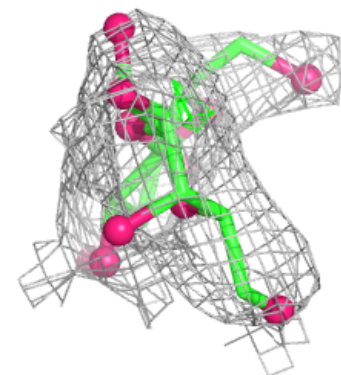
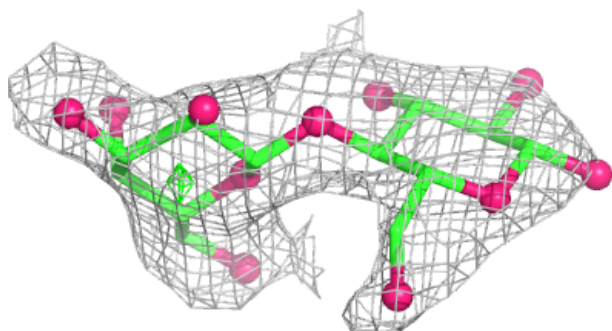
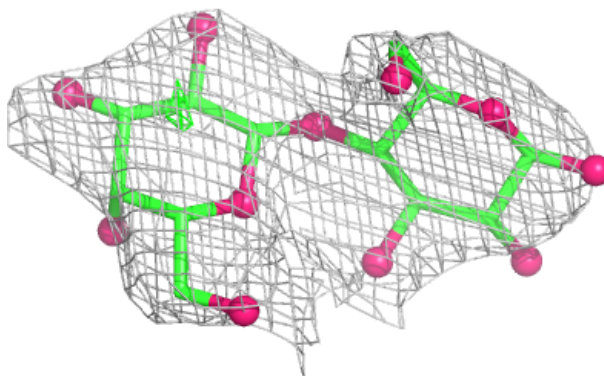


Electron density around Chain J:

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

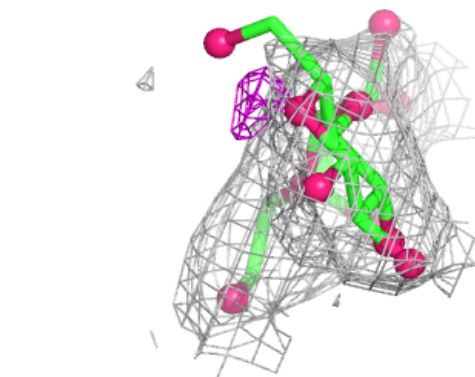
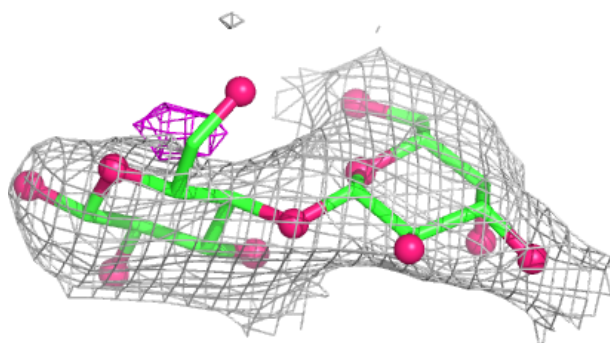
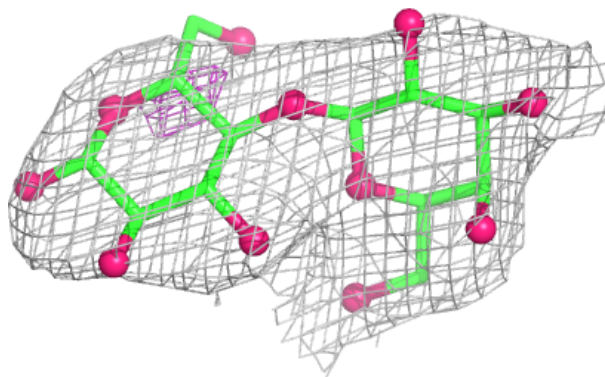
**Electron density around Chain K:**

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

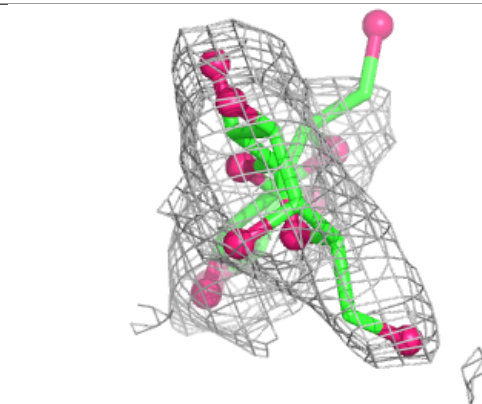
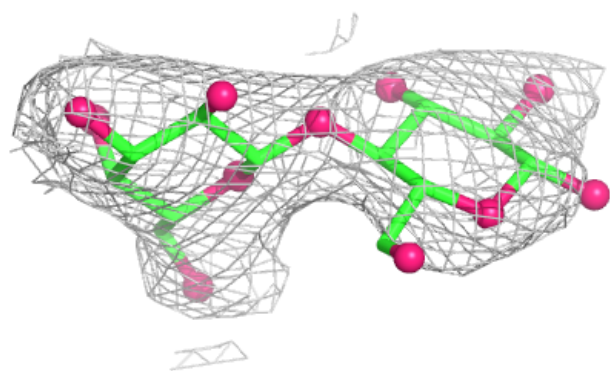
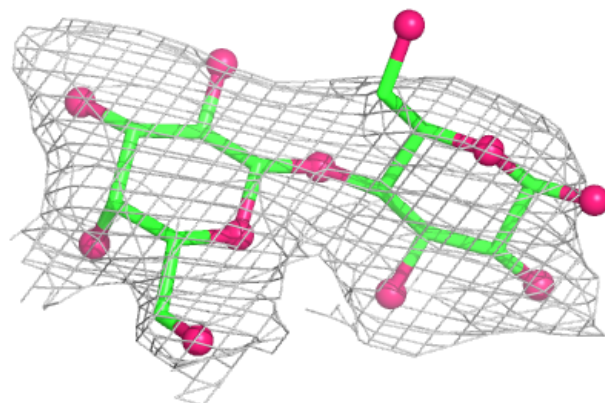


Electron density around Chain L:

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

**Electron density around Chain M:**

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



6.4 Ligands [i](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.