



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 15, 2026 – 01:23 PM UTC

PDB ID : 1L2A / pdb_0000112a
Title : The Crystal Structure and Catalytic Mechanism of Cellobiohydrolase CelS, the Major Enzymatic Component of the Clostridium thermocellum cellulosome
Authors : Guimaraes, B.G.; Souchon, H.; Lytle, B.L.; Wu, J.H.D.; Alzari, P.M.
Deposited on : 2002-02-20
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

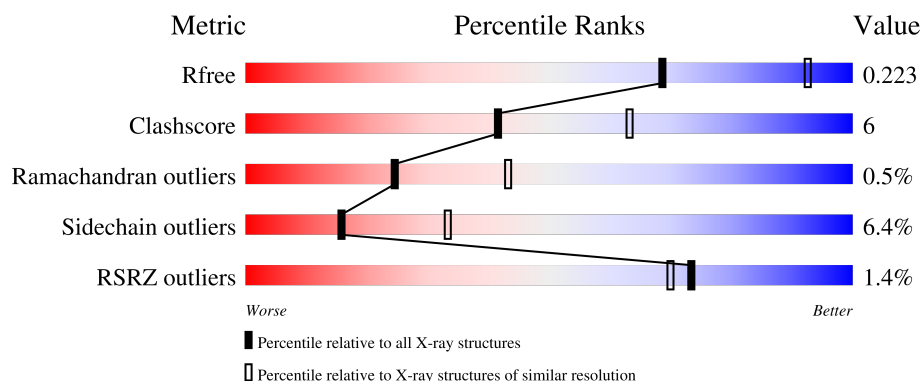
1 Overall quality at a glance

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.



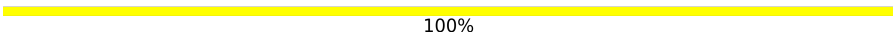
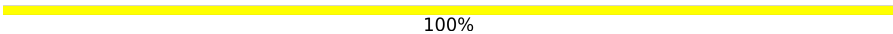
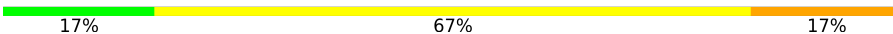
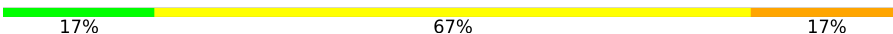
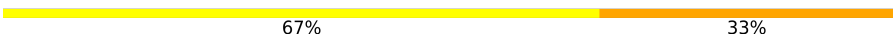
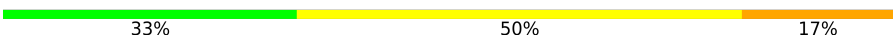
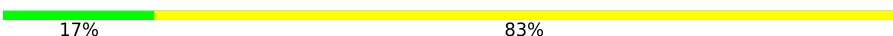
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	5829 (2.50-2.50)
Clashscore	190562	6492 (2.50-2.50)
Ramachandran outliers	187476	6378 (2.50-2.50)
Sidechain outliers	187428	6380 (2.50-2.50)
RSRZ outliers	180081	5833 (2.50-2.50)

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	678	2% 76% 16% • 5%
1	B	678	3% 77% 15% • 5%
1	C	678	% 79% 13% • 5%
1	D	678	% 82% 10% • 5%
1	E	678	% 80% 12% • 5%

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Mol	Chain	Length	Quality of chain
1	F	678	 81% 11% • 5%
2	G	2	 50% 50%
2	I	2	 50% 50%
2	K	2	 50% 50%
2	M	2	 100%
2	O	2	 100%
2	Q	2	 50% 50%
3	H	6	 17% 67% 17%
3	J	6	 17% 67% 17%
3	L	6	 67% 33%
3	N	6	 17% 33% 50%
3	P	6	 33% 50% 17%
3	R	6	 17% 83%

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 32414 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 cellobiohydrolase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	642	Total	C	N	O	S	0	0	0
			5103	3285	825	973	20			
1	B	642	Total	C	N	O	S	0	0	0
			5128	3299	835	974	20			
1	C	642	Total	C	N	O	S	0	0	0
			5109	3287	832	970	20			
1	D	642	Total	C	N	O	S	0	0	0
			5136	3303	836	977	20			
1	E	642	Total	C	N	O	S	0	0	0
			5124	3298	832	974	20			
1	F	642	Total	C	N	O	S	0	0	0
			5130	3300	833	977	20			

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



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace
2	G	2	Total	C	O	0	0	0
			22	12	10			
2	I	2	Total	C	O	0	0	0
			22	12	10			
2	K	2	Total	C	O	0	0	0
			22	12	10			
2	M	2	Total	C	O	0	0	0
			22	12	10			
2	O	2	Total	C	O	0	0	0
			22	12	10			
2	Q	2	Total	C	O	0	0	0
			22	12	10			

- Molecule 3 is an oligosaccharide called beta-D-glucopyranose-(1-4)-beta-D-glucopyranose-(1-4)-beta-D-glucopyranose-(1-4)-beta-D-glucopyranose-(1-4)-beta-D-glucopyranose.



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace
3	H	6	Total	C	O	0	0	0
			67	36	31			
3	J	6	Total	C	O	0	0	0
			67	36	31			
3	L	6	Total	C	O	0	0	0
			67	36	31			
3	N	6	Total	C	O	0	0	0
			67	36	31			
3	P	6	Total	C	O	0	0	0
			67	36	31			
3	R	6	Total	C	O	0	0	0
			67	36	31			

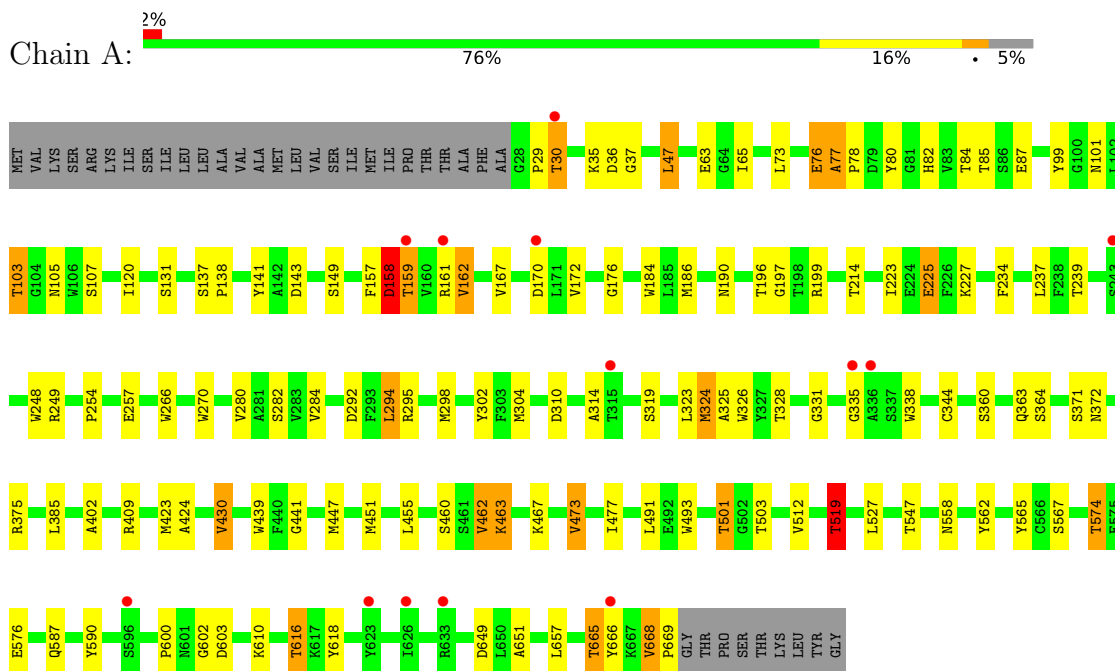
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	130	Total	O	0	0
			130	130		
4	B	127	Total	O	0	0
			127	127		
4	C	145	Total	O	0	0
			145	145		
4	D	238	Total	O	0	0
			238	238		
4	E	221	Total	O	0	0
			221	221		
4	F	289	Total	O	0	0
			289	289		

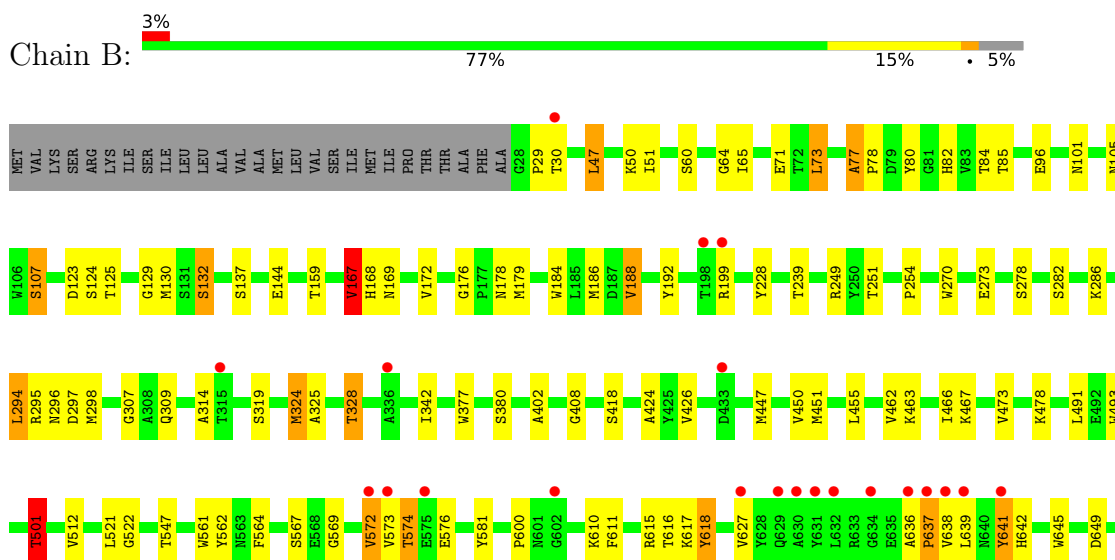
3 Residue-property plots

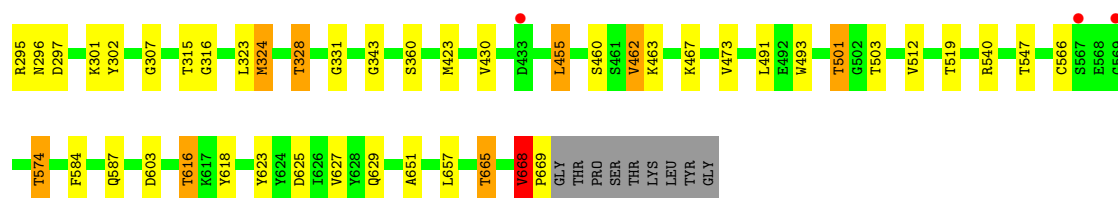
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density.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

• Molecule 1: cellobiohydrolase



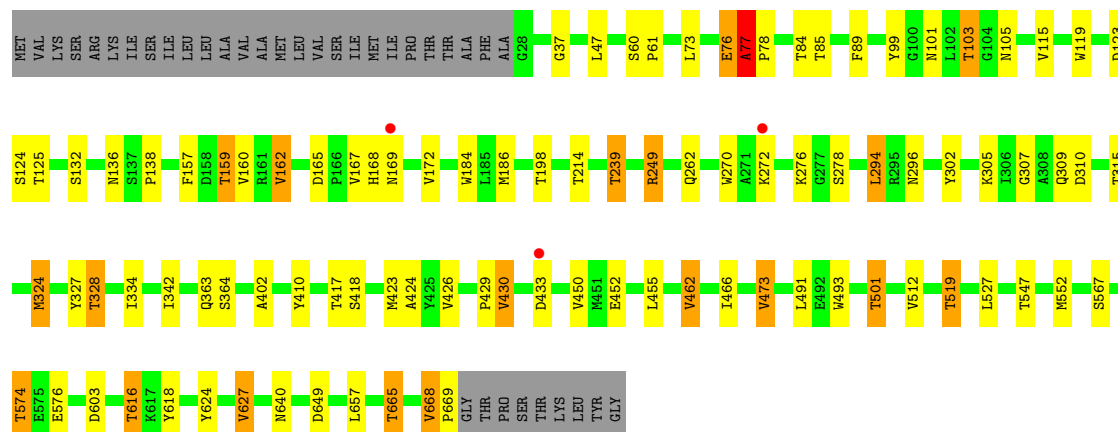
• Molecule 1: cellobiohydrolase





- Molecule 1: cellobiohydrolase

Chain F: 81% 11% 5%



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

Chain G: 50% 50%



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

Chain I: 50% 50%



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

Chain K: 50% 50%

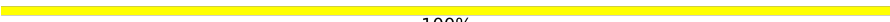


- Molecule 2: beta-D-glucopyranose-(1-4)-beta-D-glucopyranose

Chain M: 100%



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

Chain O:  100%

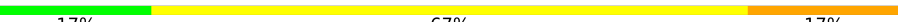
BGC1
BGC2

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

Chain Q:  50% 50%

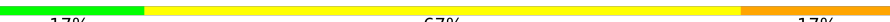
BGC1
BGC2

- Molecule 3: beta-D-glucopyranose-(1-4)-beta-D-glucopyranose-(1-4)-beta-D-glucopyranose-(1-4)-beta-D-glucopyranose-(1-4)-beta-D-glucopyranose

Chain H:  17% 67% 17%

BGC1
BGC2
BGC3
BGC4
BGC5
BGC6

- Molecule 3: beta-D-glucopyranose-(1-4)-beta-D-glucopyranose-(1-4)-beta-D-glucopyranose-(1-4)-beta-D-glucopyranose-(1-4)-beta-D-glucopyranose

Chain J:  17% 67% 17%

BGC1
BGC2
BGC3
BGC4
BGC5
BGC6

- Molecule 3: beta-D-glucopyranose-(1-4)-beta-D-glucopyranose-(1-4)-beta-D-glucopyranose-(1-4)-beta-D-glucopyranose-(1-4)-beta-D-glucopyranose

Chain L:  67% 33%

BGC1
BGC2
BGC3
BGC4
BGC5
BGC6

- Molecule 3: beta-D-glucopyranose-(1-4)-beta-D-glucopyranose-(1-4)-beta-D-glucopyranose-(1-4)-beta-D-glucopyranose-(1-4)-beta-D-glucopyranose

Chain N:  17% 33% 50%

BGC1
BGC2
BGC3
BGC4
BGC5
BGC6

- Molecule 3: beta-D-glucopyranose-(1-4)-beta-D-glucopyranose-(1-4)-beta-D-glucopyranose-(1-4)-beta-D-glucopyranose-(1-4)-beta-D-glucopyranose

Chain P:  33% 50% 17%

BGC1
BGC2
BGC3
BGC4
BGC5
BGC6

- Molecule 3: beta-D-glucopyranose-(1-4)-beta-D-glucopyranose-(1-4)-beta-D-glucopyranose-(1-4)-beta-D-glucopyranose-(1-4)-beta-D-glucopyranose-(1-4)-beta-D-glucopyranose

Chain R:  17% 83%

BGC1
BGC2
BGC3
BGC4
BGC5
BGC6

4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	148.03Å 207.64Å 215.35Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	15.00 – 2.50 15.00 – 2.50	Depositor EDS
% Data completeness (in resolution range)	95.5 (15.00-2.50) 95.0 (15.00-2.50)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.00 (at 2.51Å)	Xtriage
Refinement program	REFMAC 5.0	Depositor
R, R_{free}	0.181 , 0.226 0.181 , 0.223	Depositor DCC
R_{free} test set	10814 reflections (4.98%)	wwPDB-VP
Wilson B-factor (Å ²)	33.7	Xtriage
Anisotropy	0.085	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 28.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	0.004 for -h,l,k	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	32414	wwPDB-VP
Average B, all atoms (Å ²)	35.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 1.87% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: BGC

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	0.69	1/5282 (0.0%)	1.00	4/7213 (0.1%)
1	B	0.70	1/5307 (0.0%)	0.97	9/7241 (0.1%)
1	C	0.70	1/5288 (0.0%)	0.96	6/7218 (0.1%)
1	D	0.76	1/5315 (0.0%)	0.98	3/7251 (0.0%)
1	E	0.77	1/5303 (0.0%)	0.98	3/7236 (0.0%)
1	F	0.78	4/5309 (0.1%)	1.01	7/7244 (0.1%)
All	All	0.73	9/31804 (0.0%)	0.98	32/43403 (0.1%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	C	0	1
1	D	0	1
1	F	0	1
All	All	0	3

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	E	324	MET	SD-CE	-9.56	1.55	1.79
1	D	324	MET	SD-CE	-9.53	1.55	1.79
1	C	324	MET	SD-CE	-9.53	1.55	1.79
1	F	324	MET	SD-CE	-8.45	1.58	1.79
1	A	324	MET	SD-CE	-7.60	1.60	1.79
1	B	324	MET	SD-CE	-7.49	1.60	1.79
1	F	77	ALA	CA-C	5.33	1.59	1.52
1	F	627	VAL	CA-CB	5.21	1.61	1.54
1	F	552	MET	SD-CE	5.03	1.92	1.79

All (32) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	668	VAL	N-CA-CB	8.75	118.22	110.08
1	A	76	GLU	N-CA-C	-7.03	96.89	108.07
1	A	668	VAL	N-CA-CB	6.91	116.51	110.08
1	F	76	GLU	N-CA-C	-6.72	98.95	108.23
1	C	668	VAL	N-CA-CB	6.61	117.51	110.17
1	F	315	THR	CA-C-N	-6.60	108.48	121.41
1	F	315	THR	C-N-CA	-6.60	108.48	121.41
1	D	76	GLU	N-CA-C	-6.50	99.02	107.73
1	B	637	PRO	N-CA-C	6.45	121.33	111.15
1	D	237	LEU	N-CA-C	-5.99	105.46	112.89
1	B	188	VAL	N-CA-C	5.98	116.72	110.62
1	A	519	THR	N-CA-CB	-5.75	102.69	111.66
1	B	199	ARG	N-CA-C	5.72	118.05	108.96
1	B	167	VAL	CB-CA-C	5.64	116.02	111.06
1	C	460	SER	N-CA-C	5.63	118.14	111.33
1	F	168	HIS	N-CA-C	5.55	117.02	110.97
1	C	168	HIS	N-CA-C	5.54	117.76	111.11
1	B	641	TYR	N-CA-C	5.50	117.54	109.24
1	F	334	ILE	N-CA-C	-5.49	107.09	111.81
1	C	462	VAL	CB-CA-C	5.41	117.60	110.84
1	A	29	PRO	N-CA-C	5.41	119.68	111.19
1	B	168	HIS	N-CA-C	5.27	117.45	111.02
1	B	501	THR	N-CA-C	-5.23	105.22	113.02
1	B	65	ILE	N-CA-C	5.20	114.19	108.05
1	F	157	PHE	N-CA-C	5.18	116.74	111.14
1	C	188	VAL	N-CA-C	5.17	115.38	110.42
1	D	328	THR	OG1-CB-CG2	-5.14	99.01	109.30
1	C	157	PHE	N-CA-C	5.10	117.57	111.71
1	F	624	TYR	N-CA-C	5.07	116.62	111.14
1	B	228	TYR	N-CA-C	5.06	114.96	108.34
1	E	315	THR	CA-C-N	-5.05	111.51	121.41
1	E	315	THR	C-N-CA	-5.05	111.51	121.41

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	C	77	ALA	Peptide
1	D	316	GLY	Peptide
1	F	327	TYR	Peptide

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	5103	0	4667	86	0
1	B	5128	0	4727	69	0
1	C	5109	0	4686	61	0
1	D	5136	0	4737	54	0
1	E	5124	0	4720	64	0
1	F	5130	0	4726	75	0
2	G	22	0	19	0	0
2	I	22	0	19	0	0
2	K	22	0	19	0	0
2	M	22	0	19	0	0
2	O	22	0	19	0	0
2	Q	22	0	19	0	0
3	H	67	0	57	2	0
3	J	67	0	57	1	0
3	L	67	0	57	2	0
3	N	67	0	57	4	0
3	P	67	0	57	1	0
3	R	67	0	57	0	0
4	A	130	0	0	6	0
4	B	127	0	0	7	0
4	C	145	0	0	3	0
4	D	238	0	0	7	0
4	E	221	0	0	8	0
4	F	289	0	0	13	0
All	All	32414	0	28719	385	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 6.

All (385) 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:638:VAL:HA	4:B:798:HOH:O	1.52	1.09
1:F:310:ASP:HB2	4:F:855:HOH:O	1.53	1.06
1:D:103:THR:HG22	1:D:105:ASN:H	1.26	1.01

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:294:LEU:HG	1:C:324:MET:HE1	1.42	1.00
1:F:423:MET:SD	4:F:968:HOH:O	2.19	0.99
1:B:574:THR:HG21	1:B:576:GLU:OE1	1.63	0.99
1:A:294:LEU:HG	1:A:324:MET:HE1	1.44	0.97
1:F:294:LEU:HG	1:F:324:MET:HE1	1.48	0.94
1:A:35:LYS:HA	1:F:136:ASN:HD22	1.33	0.92
1:A:574:THR:HG21	1:A:576:GLU:OE1	1.72	0.89
1:E:103:THR:HG22	1:E:105:ASN:H	1.38	0.87
1:E:324:MET:HE3	1:E:324:MET:HA	1.56	0.87
1:F:574:THR:HG21	1:F:576:GLU:OE1	1.75	0.85
1:B:294:LEU:HG	1:B:324:MET:HE1	1.59	0.85
1:A:76:GLU:O	1:A:78:PRO:O	1.96	0.84
1:A:669:PRO:C	4:A:714:HOH:O	2.20	0.83
1:C:574:THR:HG22	1:C:641:TYR:H	1.44	0.83
1:E:501:THR:CG2	1:F:307:GLY:O	2.28	0.82
1:C:294:LEU:HG	1:C:324:MET:CE	2.10	0.81
1:A:603:ASP:CB	4:A:806:HOH:O	2.28	0.81
1:A:363:GLN:HE22	1:D:364:SER:H	1.29	0.80
1:C:99:TYR:O	1:C:103:THR:HB	1.82	0.79
1:E:307:GLY:O	1:F:501:THR:CG2	2.32	0.78
1:E:136:ASN:HB2	4:E:716:HOH:O	1.84	0.78
1:E:574:THR:HG21	4:E:774:HOH:O	1.84	0.77
1:F:37:GLY:HA2	1:F:665:THR:CG2	2.15	0.77
1:B:309:GLN:HE21	1:D:501:THR:HG21	1.49	0.76
1:D:294:LEU:HG	1:D:324:MET:HE1	1.68	0.76
1:B:501:THR:CG2	1:D:307:GLY:O	2.33	0.76
1:C:363:GLN:HE22	1:F:364:SER:H	1.33	0.75
1:B:129:GLY:O	1:B:132:SER:HB2	1.87	0.74
1:A:294:LEU:HG	1:A:324:MET:CE	2.17	0.74
1:B:328:THR:HG22	1:B:342:ILE:O	1.88	0.74
1:A:35:LYS:HA	1:F:136:ASN:ND2	2.04	0.73
1:E:278:SER:HB3	4:E:898:HOH:O	1.88	0.72
1:A:248:TRP:CZ3	1:A:331:GLY:HA2	2.23	0.72
1:A:616:THR:HB	4:A:806:HOH:O	1.89	0.71
1:C:103:THR:HG22	1:C:105:ASN:H	1.56	0.71
1:E:99:TYR:O	1:E:103:THR:HB	1.90	0.70
1:E:297:ASP:HB2	1:E:324:MET:HE1	1.74	0.70
1:B:192:TYR:C	4:B:797:HOH:O	2.35	0.70
1:F:294:LEU:HG	1:F:324:MET:CE	2.21	0.70
1:E:254:PRO:HB2	1:E:324:MET:HE2	1.74	0.69
1:B:600:PRO:HG2	4:B:797:HOH:O	1.91	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:103:THR:CG2	1:A:105:ASN:H	2.06	0.68
1:F:37:GLY:HA2	1:F:665:THR:HG21	1.75	0.68
1:B:47:LEU:O	1:B:51:ILE:HG13	1.93	0.68
1:B:50:LYS:HD3	1:B:569:GLY:O	1.93	0.67
1:C:124:SER:HB3	1:C:178:ASN:ND2	2.10	0.67
1:C:365:ASP:OD1	1:F:272:LYS:HE3	1.95	0.67
1:F:103:THR:HG22	1:F:105:ASN:H	1.61	0.66
1:D:99:TYR:O	1:D:103:THR:HB	1.95	0.66
1:B:447:MET:O	1:B:451:MET:HG2	1.96	0.66
1:D:574:THR:HG21	1:D:576:GLU:OE1	1.97	0.65
1:A:234:PHE:O	1:A:237:LEU:HB2	1.97	0.65
1:B:101:ASN:HB2	1:B:270:TRP:CE3	2.31	0.65
1:D:103:THR:HG22	1:D:105:ASN:N	2.06	0.65
1:B:573:VAL:HG23	1:B:641:TYR:O	1.96	0.65
1:B:296:ASN:HB3	1:B:328:THR:HG21	1.78	0.65
1:F:473:VAL:HG11	1:F:527:LEU:HD11	1.77	0.65
1:A:501:THR:HB	4:A:708:HOH:O	1.97	0.64
1:E:37:GLY:HA2	1:E:665:THR:CG2	2.28	0.64
1:D:519:THR:HG23	1:D:519:THR:O	1.98	0.63
1:A:63:GLU:HG3	1:A:65:ILE:HD12	1.81	0.63
1:B:77:ALA:H	1:B:78:PRO:HD2	1.63	0.63
1:C:37:GLY:HA2	1:C:665:THR:HG23	1.81	0.62
1:A:501:THR:O	1:A:501:THR:CG2	2.47	0.62
1:B:85:THR:HA	1:B:184:TRP:O	1.99	0.62
1:E:101:ASN:HB2	1:E:270:TRP:CE3	2.35	0.62
1:F:37:GLY:HA2	1:F:665:THR:HG23	1.81	0.62
1:A:36:ASP:H	1:F:136:ASN:HD21	1.46	0.62
1:C:101:ASN:HB2	1:C:270:TRP:CE3	2.34	0.62
1:C:574:THR:HG21	1:C:576:GLU:OE1	2.01	0.61
1:D:462:VAL:HG13	1:D:466:ILE:HD12	1.82	0.61
1:F:462:VAL:HG13	1:F:466:ILE:HD12	1.82	0.60
1:C:603:ASP:OD2	1:C:616:THR:HB	2.00	0.60
1:E:158:ASP:CB	4:E:893:HOH:O	2.49	0.60
1:D:76:GLU:O	1:D:77:ALA:C	2.42	0.60
1:B:84:THR:HG22	1:B:186:MET:HB3	1.84	0.60
1:A:76:GLU:O	1:A:77:ALA:C	2.45	0.60
1:E:37:GLY:HA2	1:E:665:THR:HG23	1.82	0.60
1:F:76:GLU:O	1:F:214:THR:HG21	2.02	0.59
1:A:37:GLY:HA2	1:A:665:THR:HG23	1.84	0.59
1:E:162:VAL:HG22	1:E:302:TYR:HB3	1.84	0.59
1:F:296:ASN:HB3	1:F:328:THR:HG21	1.84	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:159:THR:HG22	4:D:688:HOH:O	2.02	0.59
1:E:323:LEU:HD12	1:E:423:MET:HE1	1.84	0.59
1:A:473:VAL:HG11	1:A:527:LEU:HD11	1.84	0.59
1:F:328:THR:HG22	1:F:342:ILE:O	2.02	0.59
1:B:501:THR:HG21	1:D:309:GLN:HE21	1.67	0.58
1:E:625:ASP:O	1:E:629:GLN:HG3	2.03	0.58
1:C:128:PRO:HD2	1:C:248:TRP:NE1	2.18	0.58
1:B:616:THR:C	1:B:618:TYR:H	2.09	0.58
1:E:296:ASN:HB3	1:E:328:THR:HG21	1.84	0.58
1:A:338:TRP:CH2	3:H:5:BGC:H6C1	2.39	0.58
1:D:84:THR:HG22	1:D:186:MET:HB3	1.84	0.57
1:A:473:VAL:CG1	1:A:527:LEU:HD11	2.34	0.57
1:B:251:THR:HG21	3:J:2:BGC:H3	1.86	0.57
1:A:138:PRO:HD2	1:A:162:VAL:CG1	2.34	0.57
1:E:501:THR:HG23	1:F:307:GLY:O	2.05	0.57
1:E:77:ALA:HB3	1:E:78:PRO:CD	2.35	0.57
1:F:272:LYS:HD2	4:F:912:HOH:O	2.05	0.57
1:A:63:GLU:CG	1:A:65:ILE:HD12	2.36	0.56
1:E:199:ARG:HB3	4:E:897:HOH:O	2.05	0.56
1:B:463:LYS:O	1:B:467:LYS:HG2	2.05	0.56
1:F:76:GLU:O	1:F:77:ALA:C	2.47	0.56
1:F:101:ASN:HB2	1:F:270:TRP:CE3	2.40	0.56
1:A:447:MET:O	1:A:451:MET:HG2	2.06	0.56
1:F:249:ARG:HD2	4:F:958:HOH:O	2.05	0.56
1:F:433:ASP:C	1:F:433:ASP:OD1	2.48	0.56
1:A:254:PRO:HB2	1:A:324:MET:HE3	1.87	0.55
1:B:73:LEU:HD12	1:B:627:VAL:HG21	1.87	0.55
1:A:103:THR:HG23	1:A:105:ASN:H	1.70	0.55
1:B:616:THR:O	1:B:618:TYR:N	2.39	0.55
1:B:64:GLY:O	4:B:767:HOH:O	2.17	0.55
1:F:501:THR:HB	4:F:689:HOH:O	2.05	0.55
1:E:166:PRO:HG3	1:E:316:GLY:O	2.06	0.55
1:E:77:ALA:HB3	1:E:78:PRO:HD3	1.89	0.55
1:D:76:GLU:O	1:D:78:PRO:O	2.25	0.54
1:B:307:GLY:O	1:D:501:THR:CG2	2.55	0.54
1:B:501:THR:HG23	1:D:307:GLY:CA	2.37	0.54
1:A:460:SER:HB2	4:A:800:HOH:O	2.07	0.54
1:A:103:THR:HG22	1:A:105:ASN:H	1.72	0.54
1:A:85:THR:HA	1:A:184:TRP:O	2.08	0.54
1:B:615:ARG:O	1:B:618:TYR:HB2	2.08	0.53
1:A:77:ALA:HB3	1:A:78:PRO:CD	2.38	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:587:GLN:O	1:A:610:LYS:HA	2.08	0.53
1:B:107:SER:HA	4:B:807:HOH:O	2.08	0.53
1:B:169:ASN:HB2	4:B:727:HOH:O	2.08	0.53
1:E:307:GLY:O	1:F:501:THR:HG22	2.07	0.53
1:A:99:TYR:O	1:A:103:THR:HB	2.09	0.53
1:E:85:THR:HA	1:E:184:TRP:O	2.08	0.53
1:F:493:TRP:CZ3	1:F:512:VAL:HB	2.44	0.53
1:A:323:LEU:HD12	1:A:423:MET:HE1	1.91	0.53
1:D:101:ASN:HB2	1:D:270:TRP:CE3	2.44	0.53
1:A:172:VAL:HA	1:A:176:GLY:O	2.09	0.53
1:E:73:LEU:HD13	1:E:584:PHE:HZ	1.74	0.52
1:B:574:THR:CG2	1:B:576:GLU:OE1	2.48	0.52
1:A:47:LEU:HD13	1:A:651:ALA:HB2	1.92	0.52
1:A:364:SER:H	1:D:363:GLN:HE22	1.58	0.52
1:B:29:PRO:HB3	1:B:564:PHE:HA	1.91	0.52
1:F:450:VAL:HG12	1:F:466:ILE:HD11	1.92	0.52
1:C:462:VAL:HG13	1:C:466:ILE:HD12	1.92	0.52
1:E:225:GLU:O	1:E:226:PHE:HB2	2.10	0.52
1:A:77:ALA:HB3	1:A:78:PRO:HD2	1.92	0.52
1:B:451:MET:HE1	1:B:466:ILE:HG23	1.92	0.52
1:D:251:THR:HG21	3:N:2:BGC:H3	1.92	0.52
1:B:71:GLU:OE2	1:B:642:HIS:ND1	2.38	0.51
1:B:616:THR:C	1:B:618:TYR:N	2.67	0.51
1:C:268:ASN:HD21	1:C:272:LYS:NZ	2.08	0.51
1:E:228:TYR:O	1:E:233:GLY:HA2	2.10	0.51
1:E:254:PRO:CB	1:E:324:MET:HE2	2.41	0.51
1:C:37:GLY:HA2	1:C:665:THR:CG2	2.39	0.51
1:C:152:PRO:HG3	1:C:236:ASP:O	2.10	0.51
1:F:162:VAL:HG22	1:F:302:TYR:HB3	1.93	0.51
1:D:294:LEU:HG	1:D:324:MET:CE	2.40	0.51
1:F:249:ARG:NH2	4:F:756:HOH:O	2.43	0.51
1:B:124:SER:HB3	1:B:178:ASN:ND2	2.26	0.51
1:A:87:GLU:HG3	4:A:715:HOH:O	2.10	0.51
1:A:149:SER:HA	1:A:590:TYR:CD2	2.45	0.51
1:B:123:ASP:OD2	1:B:125:THR:HB	2.11	0.51
1:A:101:ASN:HB2	1:A:270:TRP:CE3	2.46	0.51
1:C:162:VAL:HG22	1:C:302:TYR:HB3	1.92	0.51
1:E:503:THR:HG21	4:F:966:HOH:O	2.11	0.51
1:A:473:VAL:O	1:A:477:ILE:HG13	2.11	0.51
1:D:60:SER:HB2	4:D:917:HOH:O	2.10	0.51
1:E:239:THR:HG22	4:E:873:HOH:O	2.10	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:450:VAL:HG12	1:B:466:ILE:HD11	1.93	0.50
1:F:294:LEU:CG	1:F:324:MET:HE1	2.33	0.50
1:B:562:TYR:CE1	1:B:668:VAL:HG23	2.46	0.50
1:C:144:GLU:HG2	4:C:826:HOH:O	2.10	0.50
1:C:623:TYR:O	1:C:627:VAL:HG12	2.11	0.50
1:E:603:ASP:OD1	1:E:616:THR:HB	2.11	0.50
1:A:463:LYS:O	1:A:467:LYS:HG2	2.12	0.50
1:C:364:SER:H	1:F:363:GLN:HE22	1.59	0.50
1:F:76:GLU:O	1:F:78:PRO:O	2.29	0.50
1:B:167:VAL:HG22	1:B:295:ARG:NH1	2.26	0.50
1:C:578:ARG:O	1:C:637:PRO:HD2	2.11	0.50
1:E:133:TYR:OH	1:E:138:PRO:HB3	2.12	0.50
1:D:360:SER:HB3	4:D:742:HOH:O	2.12	0.50
1:E:221:PRO:O	1:E:249:ARG:NH2	2.45	0.49
1:A:385:LEU:HD22	1:A:462:VAL:HG21	1.93	0.49
1:B:501:THR:HG22	1:D:307:GLY:O	2.11	0.49
1:D:126:GLU:HA	4:D:913:HOH:O	2.12	0.49
1:F:99:TYR:O	1:F:103:THR:HB	2.12	0.49
1:E:307:GLY:CA	1:F:501:THR:HG23	2.43	0.49
1:A:325:ALA:HB1	1:A:344:CYS:HB3	1.95	0.49
1:D:239:THR:HG22	4:D:816:HOH:O	2.12	0.49
1:D:254:PRO:HB2	1:D:324:MET:HE3	1.93	0.49
1:F:239:THR:CG2	4:F:696:HOH:O	2.59	0.49
1:A:149:SER:HA	1:A:590:TYR:CE2	2.48	0.49
1:C:669:PRO:C	4:C:706:HOH:O	2.56	0.49
1:D:37:GLY:HA2	1:D:665:THR:CG2	2.42	0.49
1:B:80:TYR:CE2	1:B:82:HIS:HB2	2.48	0.48
1:A:105:ASN:HD22	1:A:105:ASN:C	2.20	0.48
1:B:294:LEU:HB3	1:B:377:TRP:HZ2	1.78	0.48
1:C:120:ILE:HG23	1:C:182:MET:HG2	1.94	0.48
1:E:73:LEU:HD13	1:E:584:PHE:CZ	2.48	0.48
1:A:326:TRP:CD2	1:A:430:VAL:HG21	2.49	0.48
1:B:581:TYR:CE2	1:B:639:LEU:HD21	2.49	0.48
1:D:269:LYS:HE2	1:D:273:GLU:OE2	2.13	0.48
1:F:603:ASP:OD1	1:F:616:THR:HB	2.14	0.48
1:B:501:THR:HG23	1:D:307:GLY:O	2.11	0.48
1:D:519:THR:O	1:D:519:THR:CG2	2.61	0.48
1:A:501:THR:O	1:A:501:THR:HG22	2.14	0.48
4:D:767:HOH:O	3:N:1:BGC:H6C1	2.14	0.48
1:F:305:LYS:HD2	4:F:812:HOH:O	2.14	0.48
1:F:310:ASP:HB3	4:F:770:HOH:O	2.14	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:501:THR:HG22	1:F:307:GLY:O	2.12	0.48
1:B:144:GLU:HB2	1:B:408:GLY:HA3	1.95	0.48
1:F:310:ASP:HA	1:F:410:TYR:HB3	1.96	0.48
1:A:76:GLU:O	1:A:78:PRO:N	2.47	0.47
1:D:603:ASP:OD2	1:D:616:THR:HB	2.14	0.47
1:E:501:THR:HG21	1:F:309:GLN:HE21	1.78	0.47
1:A:372:ASN:HD22	1:A:375:ARG:NH2	2.12	0.47
1:E:248:TRP:CZ3	1:E:331:GLY:HA2	2.50	0.47
1:F:85:THR:HA	1:F:184:TRP:O	2.13	0.47
1:A:162:VAL:HG22	1:A:302:TYR:HB3	1.96	0.47
1:A:562:TYR:HB2	1:A:666:TYR:OH	2.14	0.47
1:C:85:THR:HA	1:C:184:TRP:O	2.14	0.47
1:B:298:MET:HE2	1:B:380:SER:CB	2.43	0.47
1:B:314:ALA:HB3	1:B:319:SER:HB3	1.97	0.47
1:C:462:VAL:HG22	1:C:465:LEU:HD23	1.96	0.47
1:C:249:ARG:HD3	1:C:249:ARG:N	2.30	0.47
3:N:1:BGC:H6C2	3:N:2:BGC:C1	2.45	0.47
1:A:257:GLU:HB2	1:A:294:LEU:HD11	1.96	0.47
1:C:63:GLU:HG3	1:C:65:ILE:HD12	1.96	0.47
1:C:251:THR:HG21	3:L:2:BGC:H3	1.97	0.47
1:D:76:GLU:O	1:D:214:THR:HG21	2.15	0.47
1:C:128:PRO:HD2	1:C:248:TRP:CD1	2.50	0.47
1:F:305:LYS:HG3	4:F:812:HOH:O	2.14	0.47
1:B:254:PRO:HG2	1:B:325:ALA:O	2.15	0.46
1:A:36:ASP:H	1:F:136:ASN:ND2	2.14	0.46
1:C:43:LEU:HD13	1:C:654:MET:HE2	1.97	0.46
1:A:63:GLU:HG3	1:A:65:ILE:CD1	2.46	0.46
1:C:134:ASN:HA	1:C:135:PRO:HD2	1.76	0.46
1:D:268:ASN:ND2	4:D:785:HOH:O	2.47	0.46
1:E:248:TRP:C	1:E:249:ARG:HG3	2.40	0.46
1:E:501:THR:HB	4:E:697:HOH:O	2.15	0.46
1:A:143:ASP:HB3	1:A:409:ARG:HG2	1.98	0.46
1:A:493:TRP:CZ3	1:A:512:VAL:HB	2.50	0.46
1:C:47:LEU:HD13	1:C:651:ALA:HB2	1.98	0.46
1:F:473:VAL:CG1	1:F:527:LEU:HD11	2.43	0.46
1:B:521:LEU:HD13	1:B:572:VAL:HG23	1.98	0.46
1:D:493:TRP:CZ3	1:D:512:VAL:HB	2.51	0.46
1:F:462:VAL:CG1	1:F:466:ILE:HD12	2.46	0.46
1:A:558:ASN:HA	1:A:666:TYR:HE1	1.81	0.46
1:B:478:LYS:HE3	4:B:803:HOH:O	2.14	0.46
1:D:455:LEU:HD23	1:D:537:ALA:HA	1.97	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:239:THR:HG21	3:H:4:BGC:H2	1.97	0.46
1:D:292:ASP:CG	1:D:368:PRO:HB3	2.40	0.46
1:B:493:TRP:CZ3	1:B:512:VAL:HB	2.51	0.46
1:D:184:TRP:CZ2	1:D:185:LEU:HD22	2.51	0.45
1:E:37:GLY:HA2	1:E:665:THR:HG21	1.97	0.45
1:E:210:GLU:HB2	1:E:430:VAL:O	2.16	0.45
1:E:463:LYS:O	1:E:467:LYS:HG2	2.16	0.45
1:B:105:ASN:ND2	1:B:107:SER:OG	2.47	0.45
1:C:426:VAL:HA	1:C:427:PRO:HD3	1.83	0.45
1:A:161:ARG:O	1:A:304:MET:HE1	2.15	0.45
1:A:602:GLY:O	1:A:603:ASP:C	2.59	0.45
1:C:126:GLU:C	1:C:128:PRO:HD3	2.41	0.45
1:D:162:VAL:HG22	1:D:302:TYR:HB3	1.99	0.45
1:A:439:TRP:CE2	1:A:441:GLY:HA3	2.51	0.45
1:D:220:HIS:HB2	1:D:221:PRO:HD2	1.98	0.45
1:A:77:ALA:CB	1:A:78:PRO:CD	2.94	0.45
1:E:297:ASP:CB	1:E:324:MET:HE1	2.44	0.45
1:F:84:THR:HG22	1:F:186:MET:HB3	1.98	0.45
1:F:89:PHE:HE2	1:F:115:VAL:HG12	1.81	0.45
1:C:73:LEU:HD13	1:C:584:PHE:HZ	1.80	0.45
1:F:519:THR:HG23	1:F:519:THR:O	2.17	0.45
1:E:324:MET:HE3	1:E:324:MET:CA	2.33	0.45
1:A:292:ASP:OD1	1:A:295:ARG:NH2	2.49	0.45
1:B:522:GLY:HA3	1:B:645:TRP:CZ3	2.52	0.45
1:E:251:THR:HG21	3:P:2:BGC:H3	1.99	0.45
1:A:372:ASN:HD22	1:A:375:ARG:HH22	1.65	0.44
1:A:519:THR:HG23	1:A:519:THR:O	2.17	0.44
1:C:166:PRO:HG3	1:C:316:GLY:O	2.18	0.44
1:C:228:TYR:O	1:C:233:GLY:HA2	2.17	0.44
1:D:328:THR:HG23	1:D:342:ILE:O	2.17	0.44
1:C:254:PRO:HB2	1:C:324:MET:HE3	1.98	0.44
1:E:668:VAL:HA	1:E:669:PRO:HD3	1.73	0.44
1:F:668:VAL:HA	1:F:669:PRO:HD3	1.80	0.44
1:A:190:ASN:HB2	1:A:199:ARG:O	2.18	0.44
1:E:84:THR:HG22	1:E:186:MET:HB3	1.98	0.44
1:B:610:LYS:O	1:B:611:PHE:C	2.58	0.44
1:C:80:TYR:CE2	1:C:82:HIS:HB2	2.53	0.44
1:C:268:ASN:HD21	1:C:272:LYS:HZ2	1.66	0.44
1:E:47:LEU:HD13	1:E:651:ALA:HB2	2.00	0.44
1:F:462:VAL:HG13	1:F:466:ILE:CD1	2.47	0.44
1:C:239:THR:HG21	3:L:4:BGC:H2	1.99	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:84:THR:HG22	1:C:186:MET:HB3	2.00	0.43
1:C:451:MET:HE1	1:C:466:ILE:HG23	2.01	0.43
1:C:668:VAL:HG23	1:C:669:PRO:HD2	2.00	0.43
1:F:159:THR:HG22	1:F:160:VAL:HG23	1.99	0.43
1:F:169:ASN:HA	1:F:172:VAL:HB	2.00	0.43
1:B:60:SER:HB3	1:B:188:VAL:HG11	1.99	0.43
1:A:402:ALA:HA	1:A:424:ALA:O	2.19	0.43
1:F:138:PRO:HD2	1:F:162:VAL:CG1	2.48	0.43
1:F:423:MET:HE2	4:F:716:HOH:O	2.18	0.43
1:A:84:THR:HG22	1:A:186:MET:HB3	1.99	0.43
1:C:519:THR:HG23	1:C:519:THR:O	2.17	0.43
1:A:314:ALA:HB3	1:A:319:SER:HB3	2.01	0.43
1:B:172:VAL:HA	1:B:176:GLY:O	2.19	0.43
1:B:254:PRO:HG3	1:B:297:ASP:OD1	2.19	0.43
1:F:402:ALA:HA	1:F:424:ALA:O	2.19	0.43
1:A:248:TRP:CE3	1:A:331:GLY:HA2	2.54	0.43
1:D:32:ALA:HB2	1:D:565:TYR:HB2	2.01	0.43
1:E:493:TRP:CZ3	1:E:512:VAL:HB	2.53	0.43
1:A:30:THR:HG21	1:A:567:SER:HB3	1.99	0.43
1:A:158:ASP:OD1	1:A:158:ASP:N	2.52	0.43
1:A:295:ARG:HA	1:A:298:MET:HE3	1.99	0.43
1:E:198:THR:CG2	4:E:902:HOH:O	2.66	0.43
1:C:165:ASP:OD1	1:C:165:ASP:C	2.61	0.43
1:D:151:TYR:HA	1:D:152:PRO:C	2.44	0.43
1:D:156:LYS:HB3	1:D:159:THR:HG23	2.00	0.43
1:A:30:THR:O	1:A:565:TYR:HB3	2.19	0.42
1:A:141:TYR:CD2	1:A:141:TYR:C	2.97	0.42
1:A:80:TYR:CE2	1:A:82:HIS:HB2	2.54	0.42
1:B:130:MET:C	1:B:132:SER:H	2.27	0.42
1:C:429:PRO:C	1:C:430:VAL:HG23	2.44	0.42
1:D:239:THR:HG21	3:N:4:BGC:H2	2.02	0.42
1:E:301:LYS:HA	1:E:343:GLY:C	2.45	0.42
1:F:574:THR:O	1:F:640:ASN:HA	2.19	0.42
1:A:37:GLY:HA2	1:A:665:THR:CG2	2.49	0.42
1:E:165:ASP:OD1	1:E:165:ASP:C	2.62	0.42
1:C:184:TRP:CE2	1:C:185:LEU:HB2	2.54	0.42
1:D:606:GLN:O	1:D:609:ILE:HG12	2.19	0.42
1:A:280:VAL:O	1:A:284:VAL:HG23	2.19	0.42
1:B:96:GLU:OE1	1:B:96:GLU:HA	2.19	0.42
1:C:363:GLN:NE2	1:F:364:SER:H	2.09	0.42
1:B:77:ALA:HB3	1:B:78:PRO:CD	2.50	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:163:GLY:HA3	1:C:303:PHE:O	2.20	0.42
1:C:668:VAL:HA	1:C:669:PRO:HD3	1.84	0.42
1:B:636:ALA:HA	1:B:637:PRO:HD3	1.93	0.42
1:C:249:ARG:NH2	4:C:784:HOH:O	2.52	0.42
1:D:309:GLN:O	1:D:412:LYS:HE3	2.20	0.42
1:E:241:ASP:OD1	1:E:242:ARG:N	2.52	0.42
1:F:165:ASP:OD1	1:F:165:ASP:C	2.63	0.42
1:F:276:LYS:C	1:F:278:SER:H	2.28	0.42
1:E:460:SER:C	1:E:462:VAL:H	2.28	0.42
1:C:45:LEU:HD23	1:C:45:LEU:HA	1.80	0.41
1:E:53:ASP:HA	1:E:54:PRO:HD2	1.95	0.41
1:E:99:TYR:CE2	1:E:103:THR:HG21	2.55	0.41
1:A:225:GLU:C	1:A:227:LYS:H	2.28	0.41
1:B:307:GLY:O	1:D:501:THR:HG23	2.20	0.41
1:E:501:THR:HG23	1:F:307:GLY:CA	2.49	0.41
1:F:119:TRP:O	1:F:186:MET:HE1	2.20	0.41
1:D:37:GLY:HA2	1:D:665:THR:HG23	2.01	0.41
1:B:402:ALA:HA	1:B:424:ALA:O	2.20	0.41
1:F:103:THR:CG2	1:F:105:ASN:H	2.32	0.41
1:A:105:ASN:ND2	1:A:107:SER:H	2.19	0.41
1:A:105:ASN:C	1:A:105:ASN:ND2	2.79	0.41
1:B:639:LEU:C	1:B:639:LEU:HD12	2.46	0.41
1:D:668:VAL:HA	1:D:669:PRO:HD3	1.81	0.41
1:F:429:PRO:C	1:F:430:VAL:HG23	2.45	0.41
1:A:266:TRP:HB3	1:A:270:TRP:CH2	2.55	0.41
1:A:266:TRP:HB3	1:A:270:TRP:CZ3	2.56	0.41
1:B:47:LEU:HG	1:B:561:TRP:CZ3	2.56	0.41
1:B:73:LEU:CD1	1:B:627:VAL:HG21	2.51	0.41
1:B:77:ALA:H	1:B:78:PRO:CD	2.33	0.41
1:C:241:ASP:HB3	1:C:243:SER:O	2.21	0.41
1:D:447:MET:HG3	1:D:469:TRP:CE2	2.56	0.41
1:E:96:GLU:HA	1:E:96:GLU:OE1	2.21	0.41
1:F:37:GLY:CA	1:F:665:THR:HG23	2.48	0.41
1:F:123:ASP:HB2	4:F:947:HOH:O	2.21	0.41
1:A:158:ASP:HB2	1:A:159:THR:H	1.73	0.41
1:B:282:SER:O	1:B:286:LYS:HG3	2.20	0.41
1:C:346:HIS:ND1	1:C:403:THR:OG1	2.52	0.41
1:D:639:LEU:HD22	1:D:641:TYR:OH	2.21	0.40
1:E:167:VAL:HG22	1:E:295:ARG:NH1	2.35	0.40
1:F:60:SER:HA	1:F:61:PRO:HD3	1.97	0.40
1:F:417:THR:HG21	1:F:424:ALA:HB1	2.02	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:141:TYR:CZ	1:C:311:LYS:HE2	2.56	0.40
1:D:615:ARG:O	1:D:618:TYR:HB2	2.21	0.40
1:E:266:TRP:HB3	1:E:270:TRP:CZ3	2.57	0.40
1:E:455:LEU:HD22	1:E:540:ARG:NH1	2.37	0.40
1:C:63:GLU:HG3	1:C:65:ILE:CD1	2.52	0.40
1:D:43:LEU:HD23	1:D:43:LEU:HA	1.90	0.40
1:E:215:TRP:HB3	1:E:587:GLN:HG2	2.04	0.40
1:C:141:TYR:CG	1:C:142:ALA:N	2.89	0.40
1:F:310:ASP:HA	1:F:410:TYR:O	2.22	0.40
1:F:574:THR:CG2	1:F:576:GLU:OE1	2.59	0.40
1:C:103:THR:HG22	1:C:105:ASN:N	2.31	0.40
1:C:493:TRP:CZ3	1:C:512:VAL:HB	2.56	0.40
1:F:262:GLN:HE21	1:F:452:GLU:CD	2.29	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	640/678 (94%)	607 (95%)	28 (4%)	5 (1%)	16	31
1	B	640/678 (94%)	605 (94%)	33 (5%)	2 (0%)	36	55
1	C	640/678 (94%)	615 (96%)	20 (3%)	5 (1%)	16	31
1	D	640/678 (94%)	622 (97%)	16 (2%)	2 (0%)	36	55
1	E	640/678 (94%)	617 (96%)	20 (3%)	3 (0%)	24	43
1	F	640/678 (94%)	621 (97%)	18 (3%)	1 (0%)	43	63
All	All	3840/4068 (94%)	3687 (96%)	135 (4%)	18 (0%)	24	43

All (18) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	77	ALA
1	A	158	ASP
1	A	335	GLY
1	C	197	GLY
1	C	335	GLY
1	D	278	SER
1	E	278	SER
1	B	617	LYS
1	C	77	ALA
1	D	77	ALA
1	E	77	ALA
1	B	77	ALA
1	C	226	PHE
1	E	197	GLY
1	F	77	ALA
1	A	600	PRO
1	A	197	GLY
1	C	430	VAL

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	512/556 (92%)	471 (92%)	41 (8%)	11	24
1	B	519/556 (93%)	488 (94%)	31 (6%)	17	36
1	C	513/556 (92%)	481 (94%)	32 (6%)	16	34
1	D	521/556 (94%)	492 (94%)	29 (6%)	19	39
1	E	518/556 (93%)	484 (93%)	34 (7%)	15	32
1	F	520/556 (94%)	487 (94%)	33 (6%)	16	34
All	All	3103/3336 (93%)	2903 (94%)	200 (6%)	16	33

All (200) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	30	THR

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Mol	Chain	Res	Type
1	A	47	LEU
1	A	73	LEU
1	A	103	THR
1	A	120	ILE
1	A	131	SER
1	A	137	SER
1	A	157	PHE
1	A	158	ASP
1	A	159	THR
1	A	162	VAL
1	A	167	VAL
1	A	170	ASP
1	A	196	THR
1	A	214	THR
1	A	223	ILE
1	A	225	GLU
1	A	249	ARG
1	A	282	SER
1	A	294	LEU
1	A	310	ASP
1	A	328	THR
1	A	360	SER
1	A	371	SER
1	A	430	VAL
1	A	455	LEU
1	A	462	VAL
1	A	463	LYS
1	A	473	VAL
1	A	491	LEU
1	A	501	THR
1	A	503	THR
1	A	519	THR
1	A	547	THR
1	A	574	THR
1	A	616	THR
1	A	618	TYR
1	A	649	ASP
1	A	657	LEU
1	A	665	THR
1	A	668	VAL
1	B	30	THR
1	B	47	LEU

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Mol	Chain	Res	Type
1	B	73	LEU
1	B	107	SER
1	B	132	SER
1	B	137	SER
1	B	159	THR
1	B	167	VAL
1	B	179	MET
1	B	239	THR
1	B	249	ARG
1	B	273	GLU
1	B	278	SER
1	B	294	LEU
1	B	328	THR
1	B	418	SER
1	B	426	VAL
1	B	455	LEU
1	B	462	VAL
1	B	473	VAL
1	B	491	LEU
1	B	501	THR
1	B	547	THR
1	B	567	SER
1	B	572	VAL
1	B	574	THR
1	B	618	TYR
1	B	649	ASP
1	B	657	LEU
1	B	665	THR
1	B	668	VAL
1	C	47	LEU
1	C	63	GLU
1	C	73	LEU
1	C	103	THR
1	C	124	SER
1	C	125	THR
1	C	159	THR
1	C	161	ARG
1	C	162	VAL
1	C	198	THR
1	C	249	ARG
1	C	294	LEU
1	C	381	TYR

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Mol	Chain	Res	Type
1	C	426	VAL
1	C	455	LEU
1	C	460	SER
1	C	462	VAL
1	C	473	VAL
1	C	491	LEU
1	C	501	THR
1	C	503	THR
1	C	513	ARG
1	C	519	THR
1	C	547	THR
1	C	574	THR
1	C	616	THR
1	C	618	TYR
1	C	627	VAL
1	C	649	ASP
1	C	657	LEU
1	C	665	THR
1	C	668	VAL
1	D	31	LYS
1	D	47	LEU
1	D	63	GLU
1	D	73	LEU
1	D	103	THR
1	D	124	SER
1	D	159	THR
1	D	162	VAL
1	D	167	VAL
1	D	198	THR
1	D	239	THR
1	D	285	SER
1	D	294	LEU
1	D	426	VAL
1	D	455	LEU
1	D	462	VAL
1	D	473	VAL
1	D	491	LEU
1	D	501	THR
1	D	519	THR
1	D	544	LYS
1	D	547	THR
1	D	574	THR

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Mol	Chain	Res	Type
1	D	616	THR
1	D	618	TYR
1	D	627	VAL
1	D	649	ASP
1	D	657	LEU
1	D	665	THR
1	E	47	LEU
1	E	63	GLU
1	E	72	THR
1	E	73	LEU
1	E	103	THR
1	E	124	SER
1	E	131	SER
1	E	159	THR
1	E	162	VAL
1	E	167	VAL
1	E	196	THR
1	E	198	THR
1	E	210	GLU
1	E	239	THR
1	E	278	SER
1	E	294	LEU
1	E	328	THR
1	E	360	SER
1	E	455	LEU
1	E	462	VAL
1	E	473	VAL
1	E	491	LEU
1	E	501	THR
1	E	519	THR
1	E	547	THR
1	E	566	CYS
1	E	574	THR
1	E	616	THR
1	E	618	TYR
1	E	623	TYR
1	E	627	VAL
1	E	657	LEU
1	E	665	THR
1	E	668	VAL
1	F	47	LEU
1	F	73	LEU

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Mol	Chain	Res	Type
1	F	103	THR
1	F	124	SER
1	F	125	THR
1	F	132	SER
1	F	159	THR
1	F	162	VAL
1	F	167	VAL
1	F	198	THR
1	F	239	THR
1	F	249	ARG
1	F	294	LEU
1	F	328	THR
1	F	418	SER
1	F	426	VAL
1	F	430	VAL
1	F	455	LEU
1	F	462	VAL
1	F	473	VAL
1	F	491	LEU
1	F	501	THR
1	F	519	THR
1	F	547	THR
1	F	567	SER
1	F	574	THR
1	F	616	THR
1	F	618	TYR
1	F	627	VAL
1	F	649	ASP
1	F	657	LEU
1	F	665	THR
1	F	668	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (44) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	101	ASN
1	A	105	ASN
1	A	207	GLN
1	A	247	GLN
1	A	363	GLN
1	A	438	GLN
1	A	640	ASN

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Mol	Chain	Res	Type
1	B	169	ASN
1	B	178	ASN
1	B	232	ASN
1	B	268	ASN
1	B	309	GLN
1	B	363	GLN
1	B	509	ASN
1	B	640	ASN
1	C	68	HIS
1	C	178	ASN
1	C	207	GLN
1	C	268	ASN
1	C	363	GLN
1	C	438	GLN
1	C	563	ASN
1	C	629	GLN
1	C	640	ASN
1	D	101	ASN
1	D	105	ASN
1	D	178	ASN
1	D	262	GLN
1	D	268	ASN
1	D	363	GLN
1	D	640	ASN
1	E	101	ASN
1	E	178	ASN
1	E	309	GLN
1	E	363	GLN
1	E	438	GLN
1	E	640	ASN
1	F	105	ASN
1	F	136	ASN
1	F	169	ASN
1	F	207	GLN
1	F	363	GLN
1	F	438	GLN
1	F	640	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 ⓘ

48 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	G	1	2	11,11,12	0.60	0	15,15,17	1.04	2 (13%)
2	BGC	G	2	2	11,11,12	0.74	0	15,15,17	0.85	0
3	BGC	H	1	3	12,12,12	0.45	0	17,17,17	1.70	5 (29%)
3	BGC	H	2	3	11,11,12	0.73	0	15,15,17	0.62	0
3	BGC	H	3	3	11,11,12	0.78	0	15,15,17	1.04	1 (6%)
3	BGC	H	4	3	11,11,12	0.58	0	15,15,17	1.78	5 (33%)
3	BGC	H	5	3	11,11,12	0.72	0	15,15,17	0.86	0
3	BGC	H	6	3	11,11,12	0.64	0	15,15,17	1.49	2 (13%)
2	BGC	I	1	2	11,11,12	0.63	0	15,15,17	1.17	2 (13%)
2	BGC	I	2	2	11,11,12	0.71	0	15,15,17	1.09	0
3	BGC	J	1	3	12,12,12	0.59	0	17,17,17	1.18	2 (11%)
3	BGC	J	2	3	11,11,12	0.88	1 (9%)	15,15,17	1.53	2 (13%)
3	BGC	J	3	3	11,11,12	0.52	0	15,15,17	0.70	1 (6%)
3	BGC	J	4	3	11,11,12	0.62	0	15,15,17	1.19	2 (13%)
3	BGC	J	5	3	11,11,12	0.76	0	15,15,17	0.72	0
3	BGC	J	6	3	11,11,12	0.58	0	15,15,17	1.73	2 (13%)
2	BGC	K	1	2	11,11,12	0.77	0	15,15,17	1.02	2 (13%)
2	BGC	K	2	2	11,11,12	0.52	0	15,15,17	0.54	0
3	BGC	L	1	3	12,12,12	0.53	0	17,17,17	1.09	1 (5%)
3	BGC	L	2	3	11,11,12	0.74	0	15,15,17	1.20	1 (6%)
3	BGC	L	3	3	11,11,12	0.69	0	15,15,17	1.17	1 (6%)
3	BGC	L	4	3	11,11,12	0.66	0	15,15,17	1.17	1 (6%)
3	BGC	L	5	3	11,11,12	0.64	0	15,15,17	0.99	1 (6%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	BGC	L	6	3	11,11,12	0.67	0	15,15,17	1.06	2 (13%)
2	BGC	M	1	2	11,11,12	0.71	0	15,15,17	1.17	2 (13%)
2	BGC	M	2	2	11,11,12	0.73	0	15,15,17	0.96	1 (6%)
3	BGC	N	1	3	12,12,12	0.67	0	17,17,17	2.19	5 (29%)
3	BGC	N	2	3	11,11,12	0.92	1 (9%)	15,15,17	1.57	2 (13%)
3	BGC	N	3	3	11,11,12	0.85	0	15,15,17	1.14	1 (6%)
3	BGC	N	4	3	11,11,12	1.00	1 (9%)	15,15,17	2.42	7 (46%)
3	BGC	N	5	3	11,11,12	0.54	0	15,15,17	0.80	0
3	BGC	N	6	3	11,11,12	0.64	0	15,15,17	1.74	1 (6%)
2	BGC	O	1	2	11,11,12	0.81	0	15,15,17	1.38	3 (20%)
2	BGC	O	2	2	11,11,12	0.87	0	15,15,17	1.10	2 (13%)
3	BGC	P	1	3	12,12,12	0.47	0	17,17,17	0.98	0
3	BGC	P	2	3	11,11,12	0.89	0	15,15,17	1.25	2 (13%)
3	BGC	P	3	3	11,11,12	0.71	0	15,15,17	0.77	0
3	BGC	P	4	3	11,11,12	0.71	0	15,15,17	1.33	4 (26%)
3	BGC	P	5	3	11,11,12	0.76	0	15,15,17	1.07	1 (6%)
3	BGC	P	6	3	11,11,12	0.68	0	15,15,17	1.40	2 (13%)
2	BGC	Q	1	2	11,11,12	0.84	0	15,15,17	1.03	2 (13%)
2	BGC	Q	2	2	11,11,12	0.52	0	15,15,17	0.58	0
3	BGC	R	1	3	12,12,12	0.69	0	17,17,17	1.32	3 (17%)
3	BGC	R	2	3	11,11,12	0.90	0	15,15,17	1.66	4 (26%)
3	BGC	R	3	3	11,11,12	0.65	0	15,15,17	1.26	1 (6%)
3	BGC	R	4	3	11,11,12	0.67	0	15,15,17	1.07	0
3	BGC	R	5	3	11,11,12	0.64	0	15,15,17	1.35	2 (13%)
3	BGC	R	6	3	11,11,12	0.66	0	15,15,17	1.48	2 (13%)

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	G	1	2	-	0/2/19/22	0/1/1/1
2	BGC	G	2	2	-	0/2/19/22	0/1/1/1
3	BGC	H	1	3	-	0/2/22/22	0/1/1/1
3	BGC	H	2	3	-	2/2/19/22	0/1/1/1
3	BGC	H	3	3	-	0/2/19/22	0/1/1/1
3	BGC	H	4	3	-	2/2/19/22	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	BGC	H	5	3	-	0/2/19/22	0/1/1/1
3	BGC	H	6	3	-	2/2/19/22	0/1/1/1
2	BGC	I	1	2	-	0/2/19/22	0/1/1/1
2	BGC	I	2	2	-	0/2/19/22	0/1/1/1
3	BGC	J	1	3	-	0/2/22/22	0/1/1/1
3	BGC	J	2	3	-	1/2/19/22	0/1/1/1
3	BGC	J	3	3	-	0/2/19/22	0/1/1/1
3	BGC	J	4	3	-	2/2/19/22	0/1/1/1
3	BGC	J	5	3	-	1/2/19/22	0/1/1/1
3	BGC	J	6	3	-	1/2/19/22	0/1/1/1
2	BGC	K	1	2	-	0/2/19/22	0/1/1/1
2	BGC	K	2	2	-	0/2/19/22	0/1/1/1
3	BGC	L	1	3	-	0/2/22/22	0/1/1/1
3	BGC	L	2	3	-	2/2/19/22	0/1/1/1
3	BGC	L	3	3	-	0/2/19/22	0/1/1/1
3	BGC	L	4	3	-	2/2/19/22	0/1/1/1
3	BGC	L	5	3	-	0/2/19/22	0/1/1/1
3	BGC	L	6	3	-	2/2/19/22	0/1/1/1
2	BGC	M	1	2	-	0/2/19/22	0/1/1/1
2	BGC	M	2	2	-	0/2/19/22	0/1/1/1
3	BGC	N	1	3	-	0/2/22/22	0/1/1/1
3	BGC	N	2	3	-	0/2/19/22	0/1/1/1
3	BGC	N	3	3	-	0/2/19/22	0/1/1/1
3	BGC	N	4	3	-	2/2/19/22	0/1/1/1
3	BGC	N	5	3	-	1/2/19/22	0/1/1/1
3	BGC	N	6	3	-	2/2/19/22	0/1/1/1
2	BGC	O	1	2	-	0/2/19/22	0/1/1/1
2	BGC	O	2	2	-	2/2/19/22	0/1/1/1
3	BGC	P	1	3	-	0/2/22/22	0/1/1/1
3	BGC	P	2	3	-	0/2/19/22	0/1/1/1
3	BGC	P	3	3	-	0/2/19/22	0/1/1/1
3	BGC	P	4	3	-	0/2/19/22	0/1/1/1
3	BGC	P	5	3	-	0/2/19/22	0/1/1/1
3	BGC	P	6	3	-	0/2/19/22	0/1/1/1
2	BGC	Q	1	2	-	0/2/19/22	0/1/1/1
2	BGC	Q	2	2	-	0/2/19/22	0/1/1/1
3	BGC	R	1	3	-	0/2/22/22	0/1/1/1
3	BGC	R	2	3	-	0/2/19/22	0/1/1/1
3	BGC	R	3	3	-	0/2/19/22	0/1/1/1
3	BGC	R	4	3	-	0/2/19/22	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	BGC	R	5	3	-	1/2/19/22	0/1/1/1
3	BGC	R	6	3	-	2/2/19/22	0/1/1/1

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	N	4	BGC	O5-C1	-2.48	1.39	1.43
3	N	2	BGC	O5-C1	-2.12	1.40	1.43
3	J	2	BGC	O5-C1	-2.03	1.40	1.43

All (82) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	N	4	BGC	C1-O5-C5	6.29	120.61	112.19
3	N	1	BGC	C6-C5-C4	-5.87	98.60	113.02
3	J	6	BGC	C1-O5-C5	5.59	119.67	112.19
3	N	6	BGC	C1-O5-C5	5.47	119.51	112.19
3	R	5	BGC	C1-O5-C5	4.37	118.04	112.19
3	N	2	BGC	O4-C4-C5	-4.14	99.12	109.32
3	R	2	BGC	C1-O5-C5	4.12	117.70	112.19
3	N	1	BGC	C3-C4-C5	3.91	117.33	110.23
3	H	1	BGC	O5-C5-C4	3.76	116.47	109.70
3	H	6	BGC	C1-O5-C5	3.75	117.21	112.19
3	P	6	BGC	C1-O5-C5	3.66	117.10	112.19
3	R	3	BGC	C1-O5-C5	3.65	117.08	112.19
3	R	6	BGC	C1-O5-C5	3.57	116.97	112.19
2	O	1	BGC	O5-C1-C2	-3.45	102.56	110.79
3	N	1	BGC	C1-O5-C5	3.40	120.23	113.65
3	J	2	BGC	C1-O5-C5	3.39	116.73	112.19
3	N	2	BGC	C1-O5-C5	3.36	116.69	112.19
3	N	4	BGC	O6-C6-C5	-3.16	100.57	111.33
3	H	4	BGC	C2-C3-C4	3.15	116.41	110.86
3	H	1	BGC	C3-C4-C5	3.15	115.94	110.23
3	N	4	BGC	C6-C5-C4	-3.08	105.44	113.02
3	H	6	BGC	C1-C2-C3	3.08	114.13	109.64
3	L	3	BGC	C1-O5-C5	3.06	116.29	112.19
3	H	3	BGC	C1-O5-C5	3.04	116.26	112.19
3	H	4	BGC	C1-O5-C5	2.96	116.16	112.19
3	N	4	BGC	O4-C4-C5	-2.84	102.33	109.32
2	I	1	BGC	O5-C1-C2	-2.84	104.02	110.79
3	N	4	BGC	O5-C5-C4	2.83	117.70	110.83
3	J	1	BGC	O5-C5-C4	2.81	114.76	109.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	O	2	BGC	O5-C1-C2	-2.73	104.28	110.79
2	Q	1	BGC	O5-C1-C2	-2.70	104.34	110.79
3	L	2	BGC	O4-C4-C5	-2.70	102.68	109.32
3	R	1	BGC	O4-C4-C3	-2.69	104.03	110.38
3	L	1	BGC	O4-C4-C3	-2.65	104.13	110.38
3	H	4	BGC	C1-C2-C3	2.64	113.49	109.64
3	H	1	BGC	C1-O5-C5	2.62	118.72	113.65
3	H	1	BGC	C6-C5-C4	-2.60	106.63	113.02
2	M	1	BGC	O5-C1-C2	-2.58	104.64	110.79
2	G	1	BGC	C1-O5-C5	2.56	115.62	112.19
3	N	1	BGC	O5-C5-C4	2.55	114.30	109.70
3	H	4	BGC	C3-C4-C5	2.52	114.80	110.23
3	J	1	BGC	C1-O5-C5	2.48	118.46	113.65
2	M	2	BGC	C1-O5-C5	2.48	115.52	112.19
3	P	2	BGC	O3-C3-C4	2.48	116.22	110.38
3	R	2	BGC	O4-C4-C5	-2.47	103.25	109.32
3	R	6	BGC	C1-C2-C3	2.46	113.23	109.64
2	O	1	BGC	C3-C4-C5	-2.46	105.77	110.23
3	N	1	BGC	O3-C3-C4	-2.44	104.62	110.38
3	J	6	BGC	C3-C4-C5	2.43	114.64	110.23
3	J	2	BGC	O4-C4-C5	-2.43	103.34	109.32
3	P	4	BGC	O3-C3-C2	-2.41	105.14	110.05
2	O	2	BGC	C1-O5-C5	2.40	115.40	112.19
3	J	4	BGC	O3-C3-C2	-2.39	105.17	110.05
3	P	4	BGC	O4-C4-C5	-2.37	103.48	109.32
3	P	4	BGC	O5-C5-C6	2.36	112.26	107.66
3	P	6	BGC	C1-C2-C3	2.33	113.04	109.64
3	R	1	BGC	O5-C1-C2	-2.32	106.22	110.30
2	M	1	BGC	C1-O5-C5	2.31	115.28	112.19
2	Q	1	BGC	O5-C5-C6	2.30	112.14	107.66
3	P	2	BGC	O4-C4-C5	-2.30	103.67	109.32
3	R	1	BGC	O5-C5-C6	2.29	112.11	106.44
3	R	2	BGC	O3-C3-C4	2.28	115.75	110.38
3	L	5	BGC	O3-C3-C2	-2.28	105.40	110.05
2	K	1	BGC	O5-C1-C2	-2.28	105.36	110.79
3	H	1	BGC	O3-C3-C4	-2.24	105.11	110.38
3	R	2	BGC	O2-C2-C3	-2.22	105.55	110.15
3	N	4	BGC	O5-C5-C6	-2.21	103.37	107.66
3	P	5	BGC	C1-O5-C5	2.16	115.08	112.19
3	J	4	BGC	O5-C5-C6	2.15	111.84	107.66
3	L	4	BGC	O3-C3-C2	-2.13	105.70	110.05
2	G	1	BGC	O5-C1-C2	-2.13	105.71	110.79

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	K	1	BGC	O5-C5-C6	2.11	111.77	107.66
3	H	4	BGC	C6-C5-C4	-2.10	107.85	113.02
3	N	4	BGC	C3-C4-C5	2.10	114.05	110.23
3	J	3	BGC	O4-C4-C3	-2.10	105.44	110.38
3	N	3	BGC	O4-C4-C5	-2.08	104.19	109.32
3	R	5	BGC	O4-C4-C5	-2.06	104.24	109.32
3	P	4	BGC	C1-O5-C5	2.06	114.94	112.19
3	L	6	BGC	C1-C2-C3	2.05	112.64	109.64
2	I	1	BGC	C1-O5-C5	2.05	114.94	112.19
3	L	6	BGC	C1-O5-C5	2.05	114.93	112.19
2	O	1	BGC	O2-C2-C3	2.01	114.31	110.15

There are no chirality outliers.

All (27) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	L	2	BGC	O5-C5-C6-O6
3	R	6	BGC	O5-C5-C6-O6
3	L	2	BGC	C4-C5-C6-O6
3	N	4	BGC	O5-C5-C6-O6
3	L	4	BGC	C4-C5-C6-O6
3	L	4	BGC	O5-C5-C6-O6
3	N	4	BGC	C4-C5-C6-O6
3	R	6	BGC	C4-C5-C6-O6
3	H	6	BGC	O5-C5-C6-O6
3	N	6	BGC	C4-C5-C6-O6
3	H	2	BGC	O5-C5-C6-O6
3	H	4	BGC	O5-C5-C6-O6
3	H	4	BGC	C4-C5-C6-O6
3	H	6	BGC	C4-C5-C6-O6
3	N	6	BGC	O5-C5-C6-O6
3	L	6	BGC	O5-C5-C6-O6
3	L	6	BGC	C4-C5-C6-O6
3	J	4	BGC	C4-C5-C6-O6
3	J	5	BGC	O5-C5-C6-O6
3	J	2	BGC	C4-C5-C6-O6
3	J	6	BGC	C4-C5-C6-O6
3	J	4	BGC	O5-C5-C6-O6
2	O	2	BGC	C4-C5-C6-O6
3	H	2	BGC	C4-C5-C6-O6
3	N	5	BGC	O5-C5-C6-O6
3	R	5	BGC	O5-C5-C6-O6

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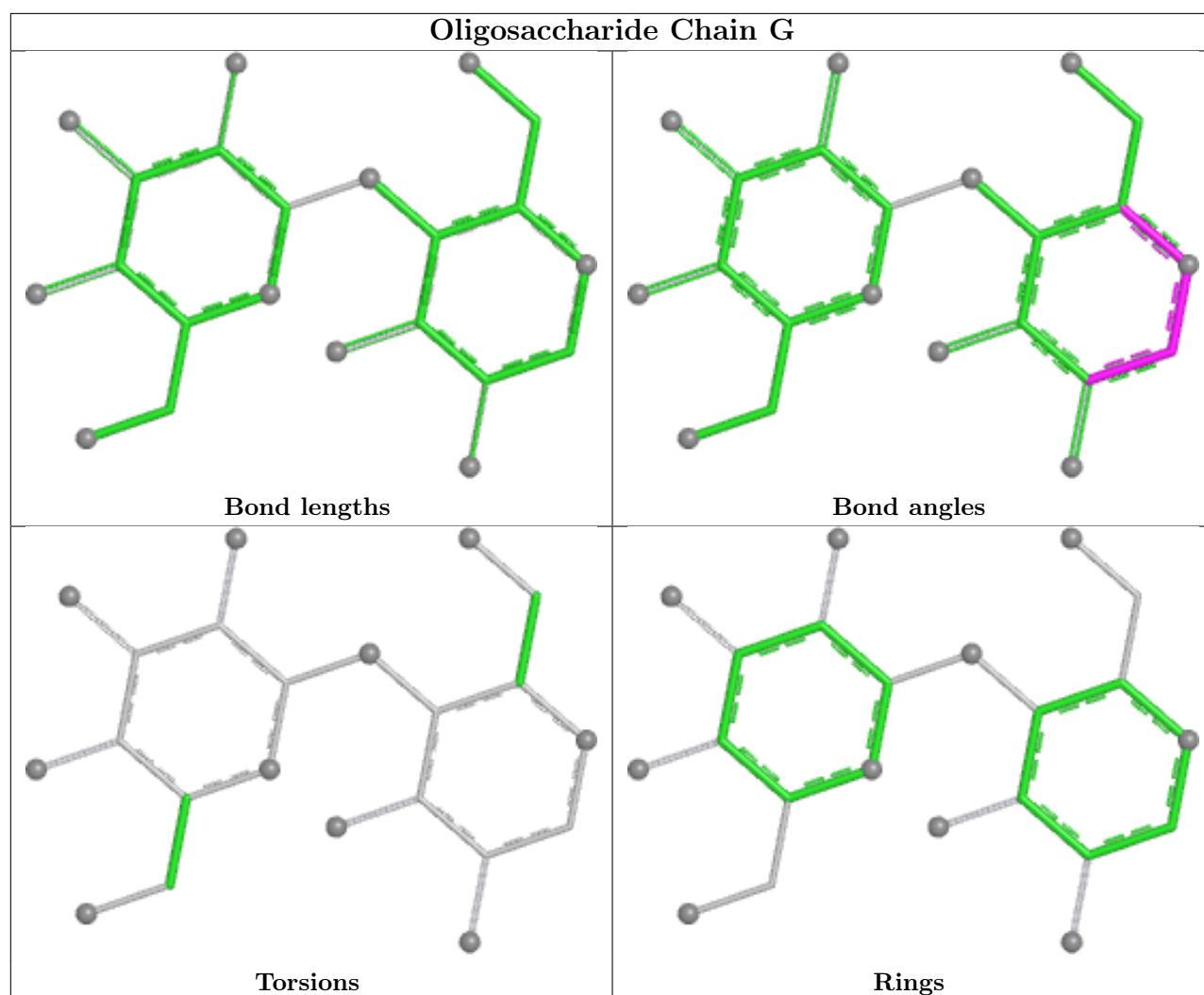
Mol	Chain	Res	Type	Atoms
2	O	2	BGC	O5-C5-C6-O6

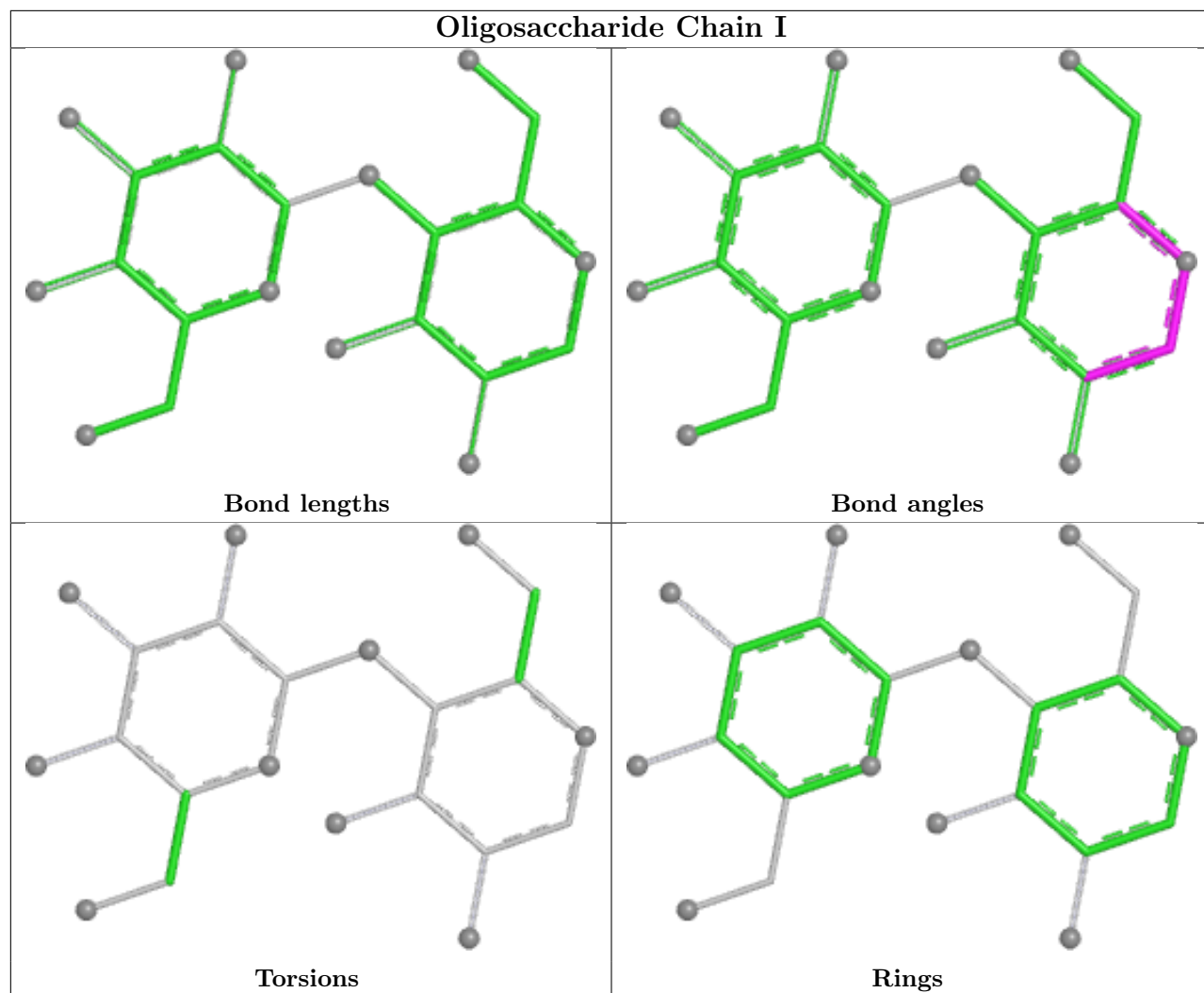
There are no ring outliers.

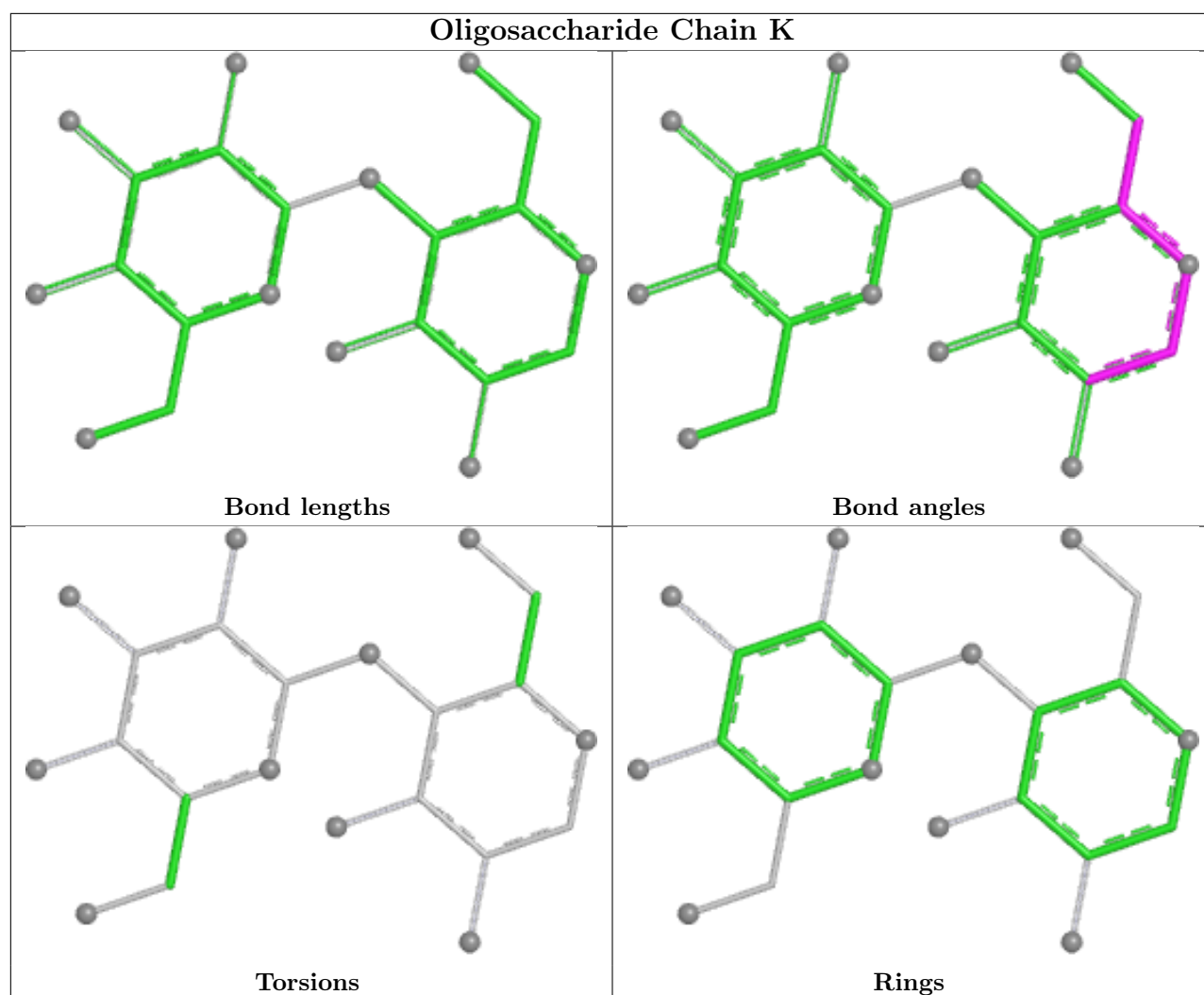
9 monomers are involved in 10 short contacts:

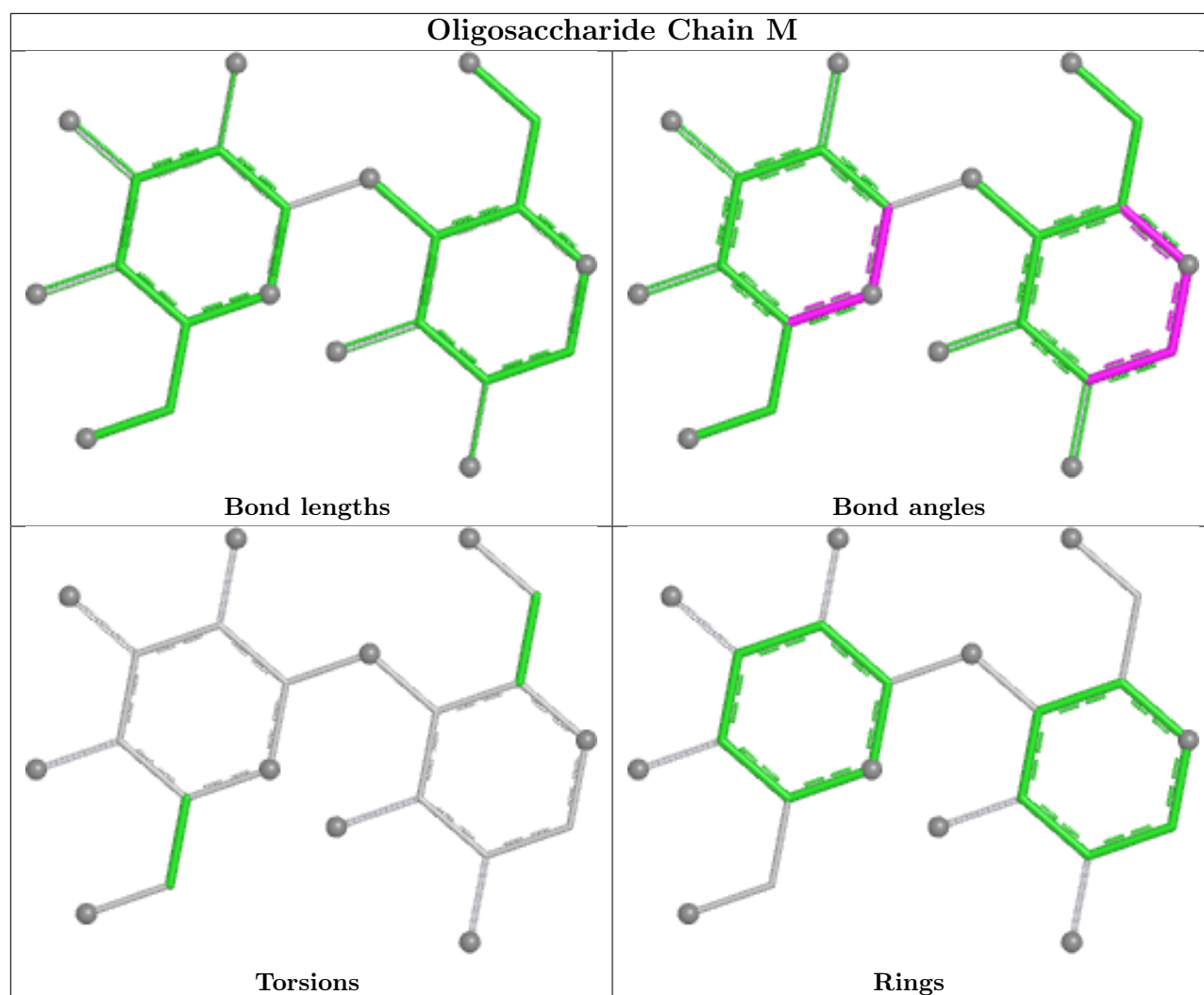
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	L	2	BGC	1	0
3	H	5	BGC	1	0
3	J	2	BGC	1	0
3	N	4	BGC	1	0
3	N	1	BGC	2	0
3	P	2	BGC	1	0
3	L	4	BGC	1	0
3	H	4	BGC	1	0
3	N	2	BGC	2	0

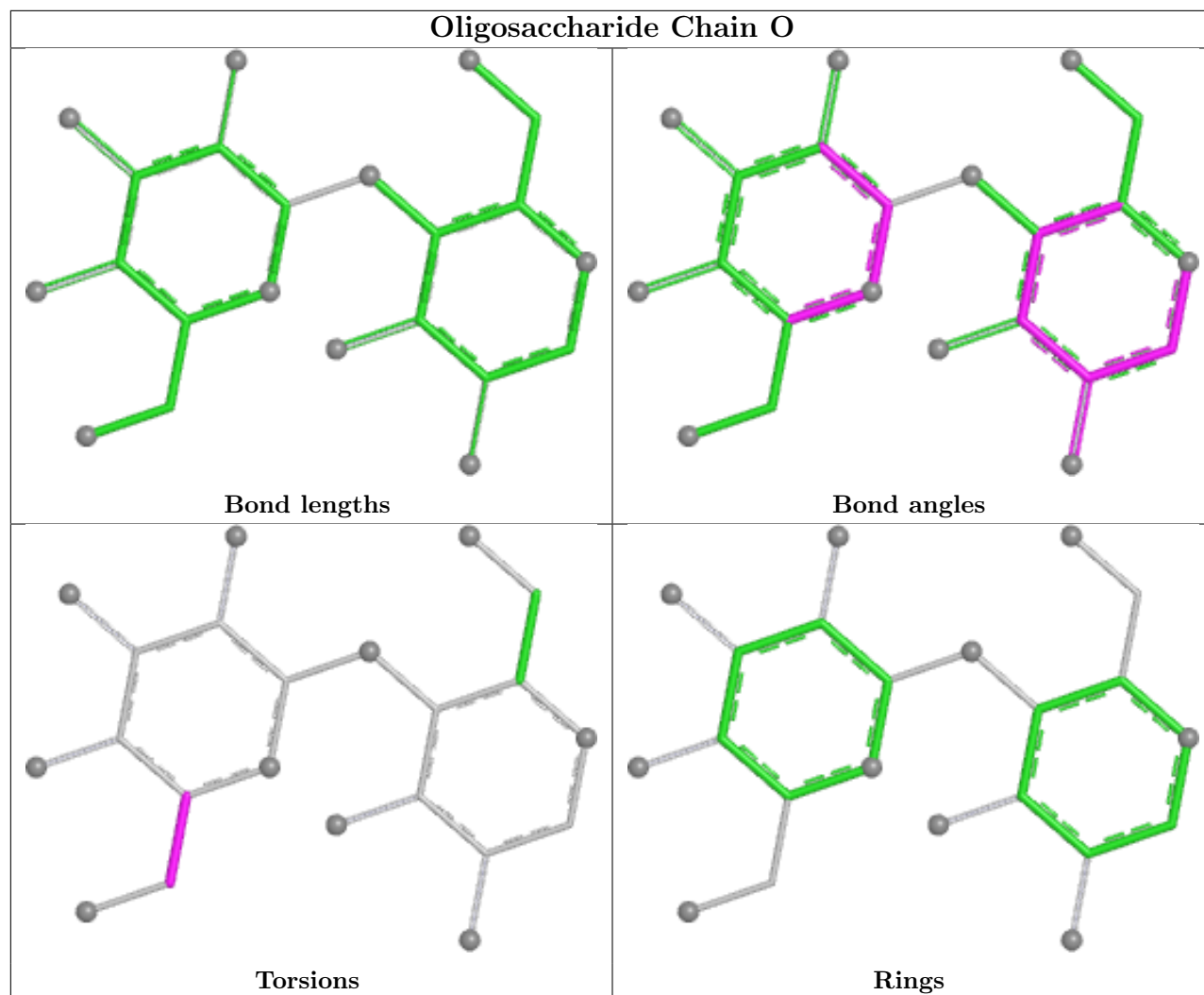
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for oligosaccharide.



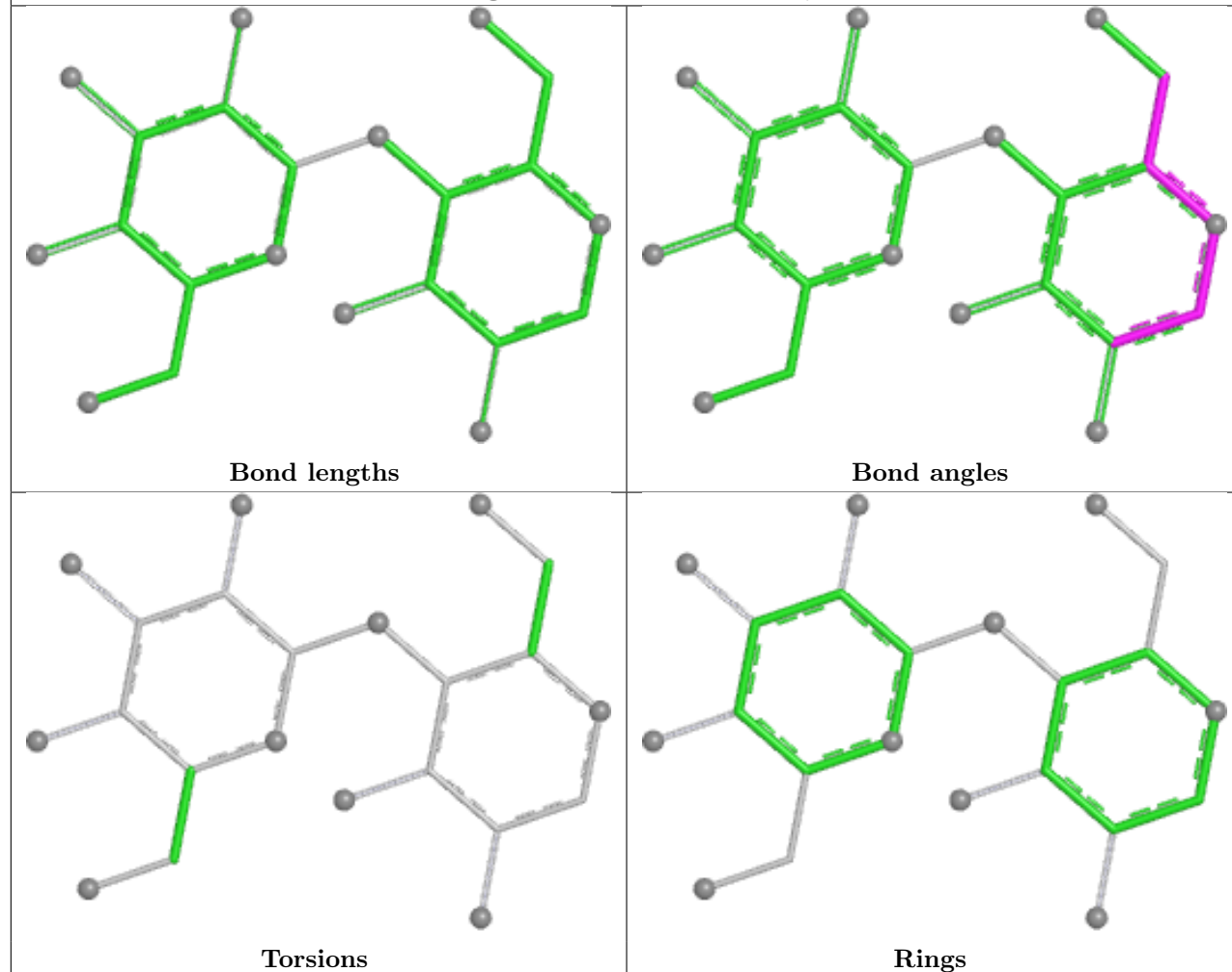




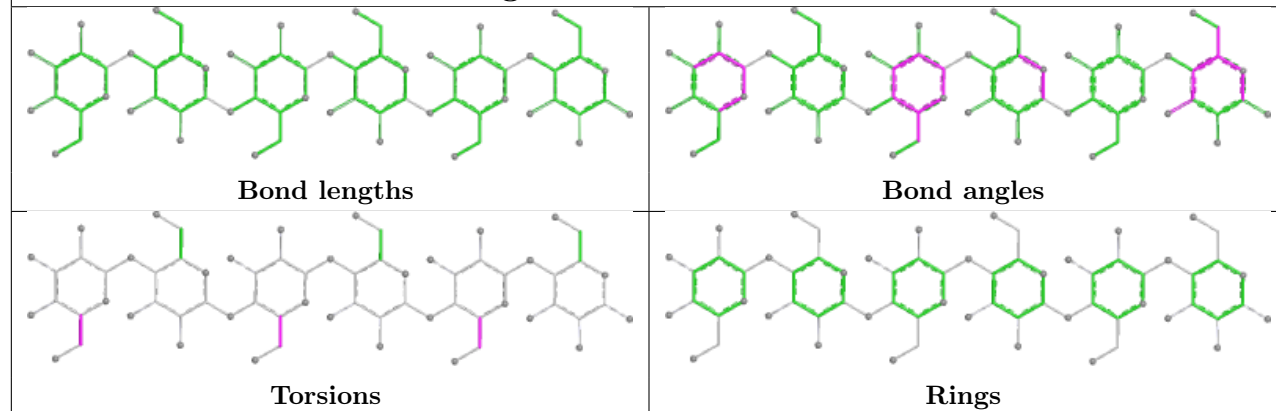


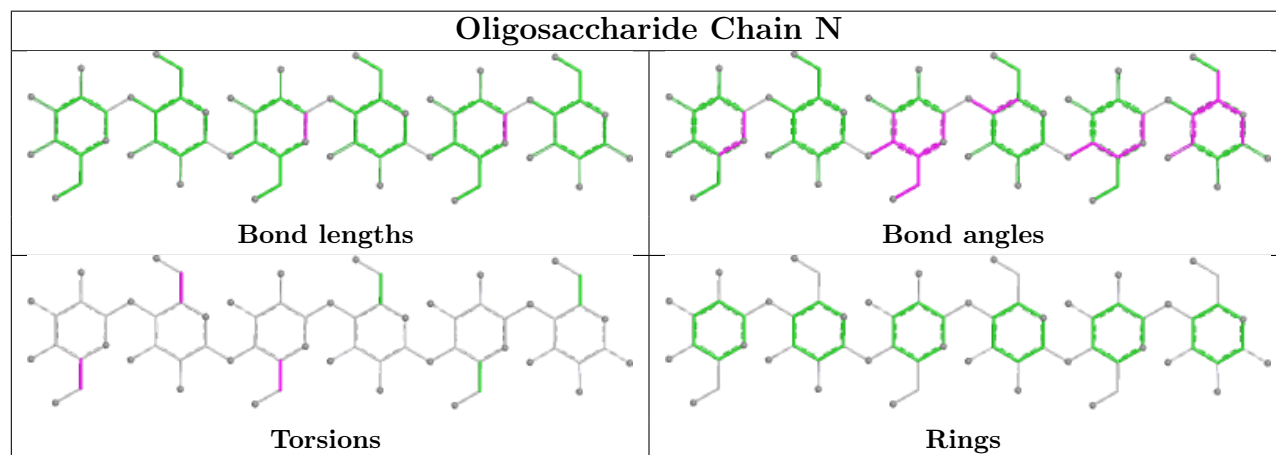
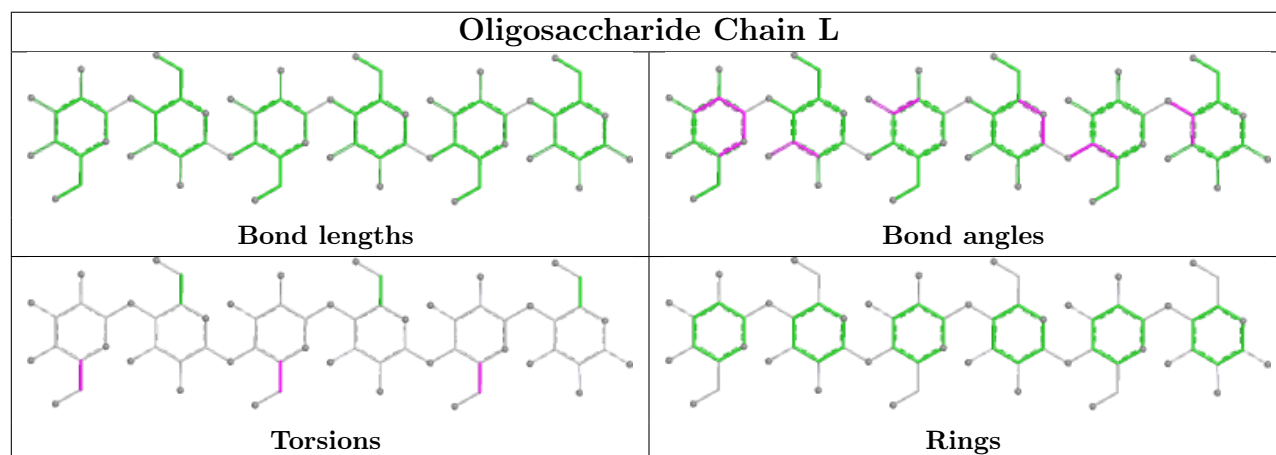
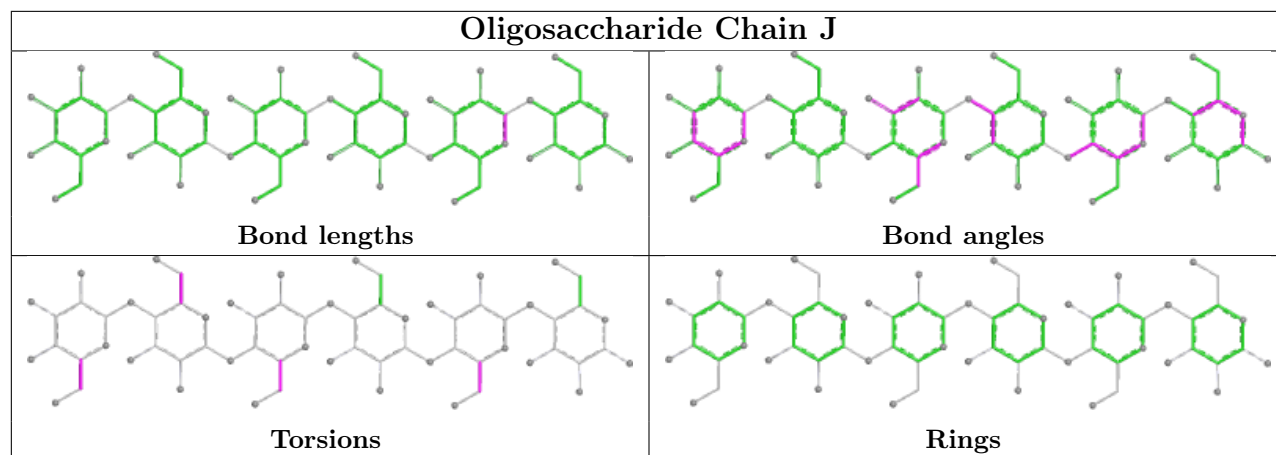


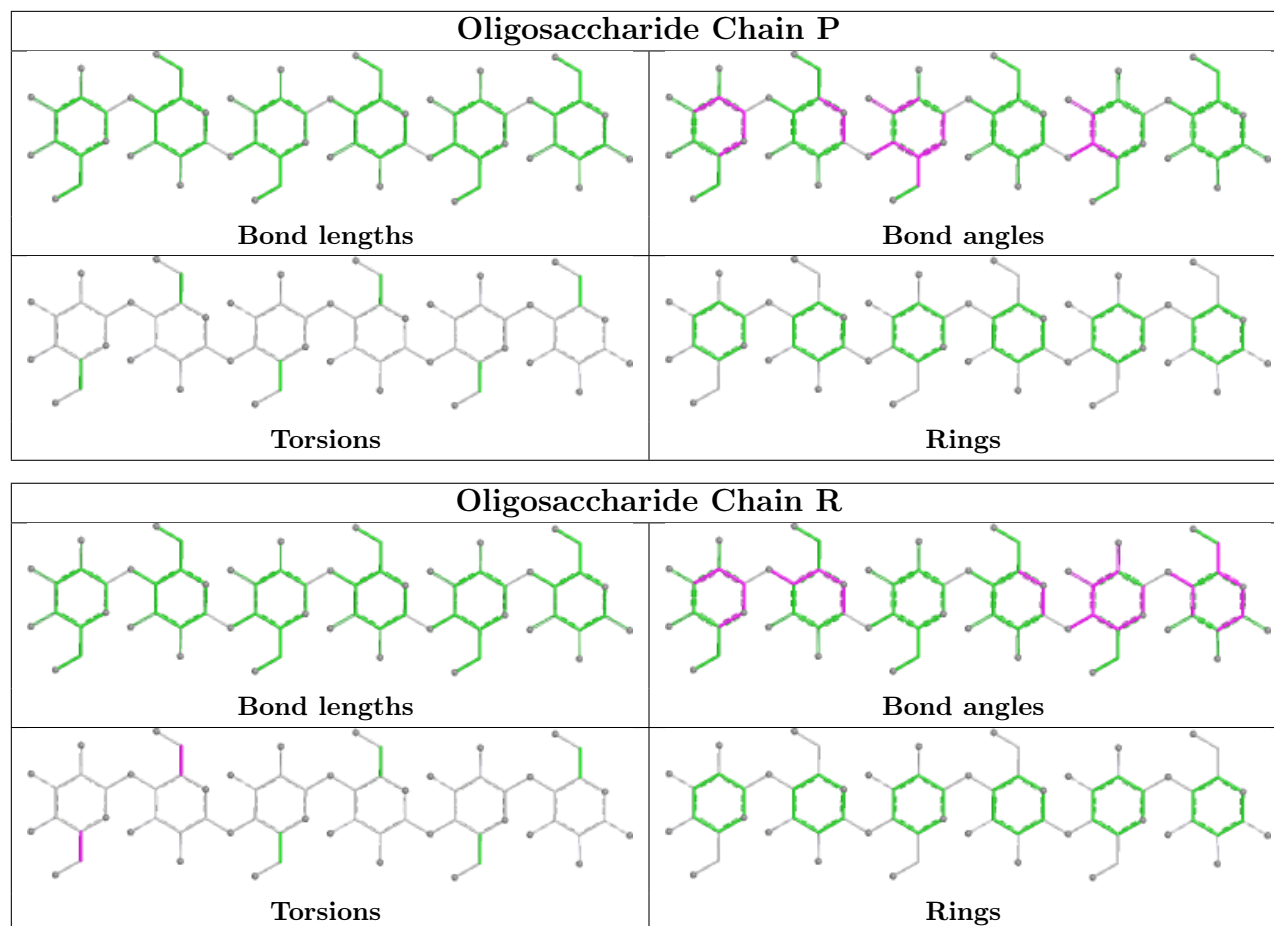
Oligosaccharide Chain Q



Oligosaccharide Chain H







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 ⓘ

6.1 Protein, DNA and RNA chains ⓘ

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	642/678 (94%)	-0.16	13 (2%) 65 61	20, 39, 62, 68	0
1	B	642/678 (94%)	-0.07	21 (3%) 49 45	23, 41, 61, 79	0
1	C	642/678 (94%)	-0.33	7 (1%) 78 75	21, 37, 61, 74	0
1	D	642/678 (94%)	-0.59	4 (0%) 85 83	20, 30, 45, 55	0
1	E	642/678 (94%)	-0.64	5 (0%) 82 80	18, 29, 42, 52	0
1	F	642/678 (94%)	-0.74	3 (0%) 87 85	17, 26, 38, 51	0
All	All	3852/4068 (94%)	-0.42	53 (1%) 73 70	17, 32, 58, 79	0

All (53) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	638	VAL	7.6
1	B	639	LEU	5.8
1	B	630	ALA	3.9
1	E	567	SER	3.6
1	B	602	GLY	3.5
1	B	433	ASP	3.2
1	D	28	GLY	3.0
1	B	573	VAL	3.0
1	A	315	THR	3.0
1	C	159	THR	3.0
1	B	198	THR	2.9
1	D	277	GLY	2.9
1	A	159	THR	2.9
1	C	334	ILE	2.9
1	B	632	LEU	2.8
1	A	30	THR	2.8
1	E	433	ASP	2.8
1	A	170	ASP	2.8
1	F	433	ASP	2.7

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Mol	Chain	Res	Type	RSRZ
1	E	277	GLY	2.7
1	B	336	ALA	2.6
1	D	316	GLY	2.6
1	A	623	TYR	2.5
1	A	626	ILE	2.5
1	B	315	THR	2.5
1	A	336	ALA	2.5
1	B	634	GLY	2.5
1	E	169	ASN	2.5
1	A	666	TYR	2.4
1	B	572	VAL	2.4
1	A	161	ARG	2.4
1	A	633	ARG	2.3
1	B	199	ARG	2.3
1	A	243	SER	2.3
1	B	636	ALA	2.3
1	F	272	LYS	2.2
1	C	158	ASP	2.2
1	B	637	PRO	2.2
1	B	631	TYR	2.2
1	B	30	THR	2.2
1	B	627	VAL	2.2
1	C	28	GLY	2.2
1	A	596	SER	2.2
1	B	575	GLU	2.2
1	B	629	GLN	2.1
1	B	641	TYR	2.1
1	C	337	SER	2.1
1	D	433	ASP	2.1
1	A	335	GLY	2.1
1	E	569	GLY	2.1
1	C	243	SER	2.1
1	F	169	ASN	2.0
1	C	336	ALA	2.0

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

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

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.

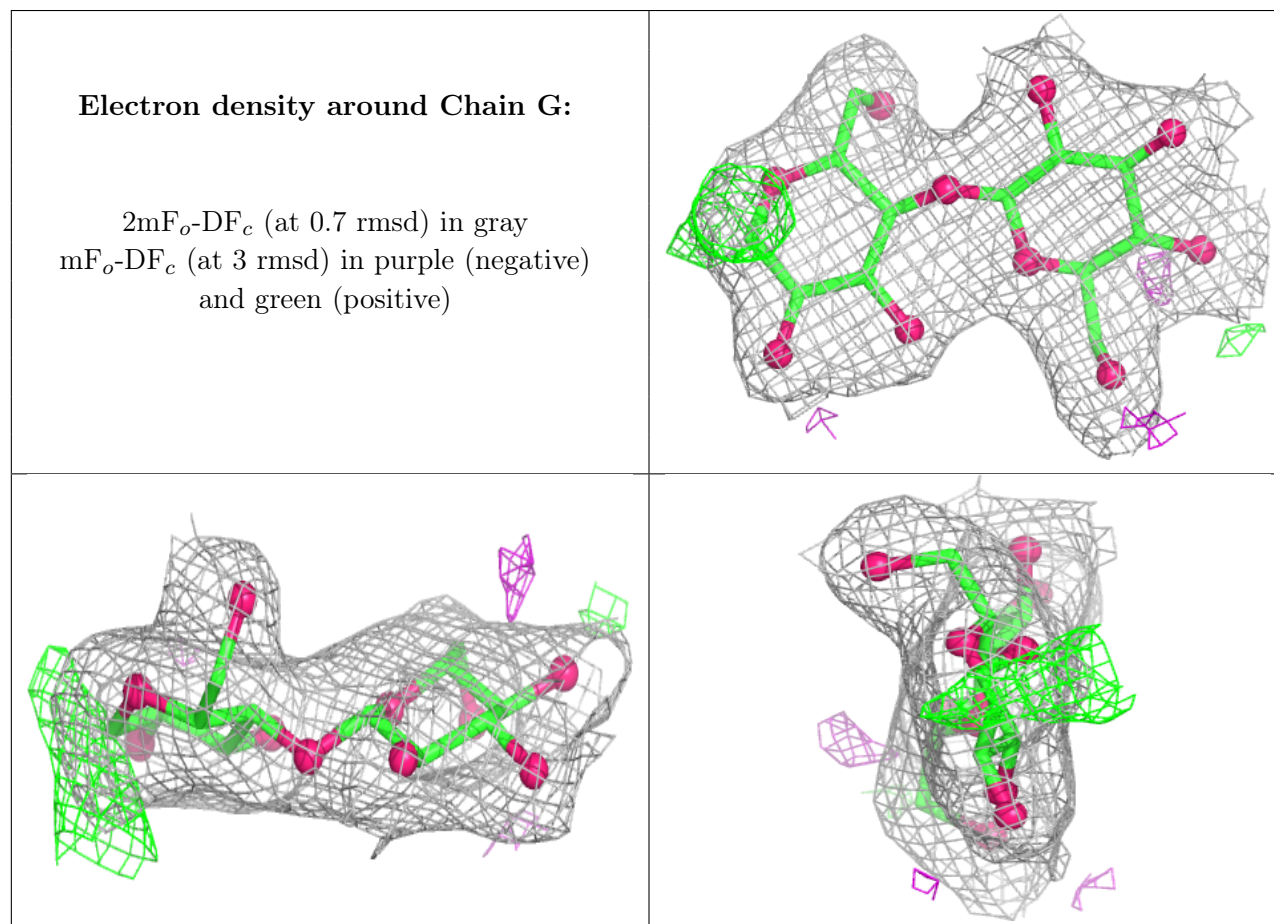
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
3	BGC	H	5	11/12	0.77	0.14	67,68,71,74	0
3	BGC	L	6	11/12	0.77	0.13	75,77,78,78	0
3	BGC	H	6	11/12	0.80	0.12	76,77,78,79	0
3	BGC	J	6	11/12	0.83	0.12	64,66,67,69	0
3	BGC	N	6	11/12	0.88	0.09	40,44,46,47	0
2	BGC	I	1	11/12	0.89	0.08	42,44,45,46	0
3	BGC	L	5	11/12	0.90	0.10	61,64,67,72	0
2	BGC	K	1	11/12	0.90	0.07	34,36,37,38	0
2	BGC	G	1	11/12	0.90	0.07	33,36,38,40	0
2	BGC	O	1	11/12	0.91	0.06	27,30,32,32	0
3	BGC	H	4	11/12	0.91	0.08	56,58,61,64	0
3	BGC	J	4	11/12	0.92	0.08	39,42,46,51	0
3	BGC	J	5	11/12	0.92	0.09	53,54,56,60	0
3	BGC	P	6	11/12	0.92	0.07	45,48,49,50	0
3	BGC	R	6	11/12	0.92	0.08	43,46,52,57	0
3	BGC	L	2	11/12	0.93	0.08	39,42,44,44	0
3	BGC	L	3	11/12	0.93	0.07	43,44,45,48	0
3	BGC	L	4	11/12	0.93	0.07	45,51,54,59	0
2	BGC	M	1	11/12	0.93	0.07	32,35,37,38	0
2	BGC	I	2	11/12	0.94	0.06	34,40,41,41	0
2	BGC	Q	1	11/12	0.94	0.07	25,27,31,32	0
3	BGC	N	1	12/12	0.94	0.07	29,30,33,40	0
3	BGC	N	4	11/12	0.94	0.07	28,31,33,36	0
3	BGC	H	1	12/12	0.94	0.07	38,42,44,47	0
3	BGC	J	3	11/12	0.94	0.07	35,37,40,41	0
3	BGC	H	3	11/12	0.94	0.07	43,47,50,53	0
3	BGC	H	2	11/12	0.95	0.07	40,43,44,45	0
3	BGC	N	3	11/12	0.96	0.06	26,28,31,32	0
3	BGC	R	2	11/12	0.96	0.06	21,26,27,27	0
3	BGC	R	3	11/12	0.96	0.05	26,27,29,29	0
3	BGC	R	4	11/12	0.96	0.06	28,29,31,31	0
3	BGC	P	5	11/12	0.96	0.06	32,35,37,42	0
3	BGC	P	3	11/12	0.97	0.04	23,28,29,29	0
3	BGC	P	4	11/12	0.97	0.04	25,29,30,31	0
3	BGC	L	1	12/12	0.97	0.05	33,38,39,40	0
2	BGC	G	2	11/12	0.97	0.05	30,31,32,32	0
3	BGC	R	1	12/12	0.97	0.05	23,25,28,30	0

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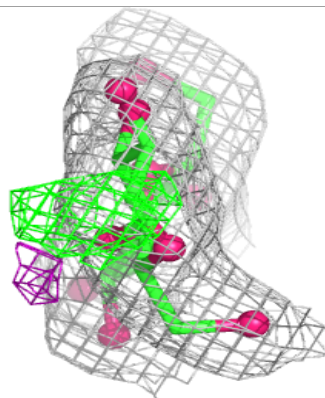
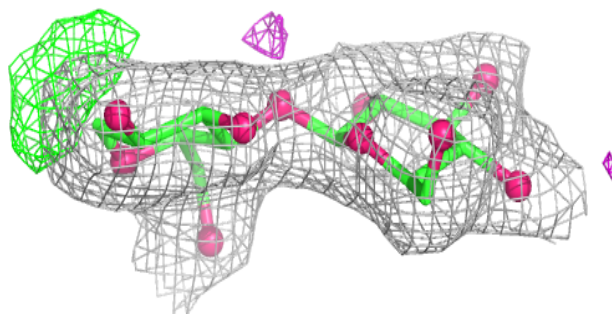
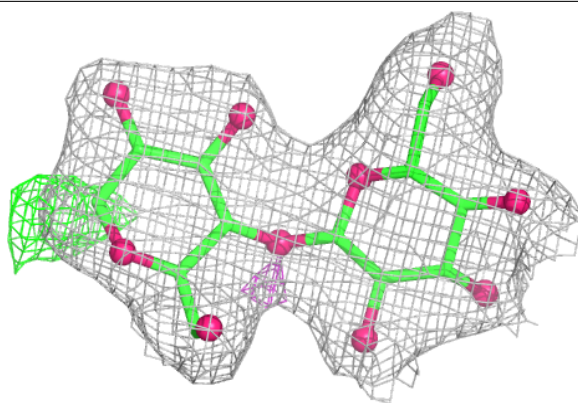
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	BGC	M	2	11/12	0.97	0.04	28,30,31,32	0
3	BGC	J	1	12/12	0.97	0.05	33,35,37,38	0
3	BGC	N	5	11/12	0.97	0.05	33,34,38,38	0
3	BGC	R	5	11/12	0.97	0.05	32,35,37,40	0
3	BGC	J	2	11/12	0.97	0.06	29,35,39,42	0
2	BGC	O	2	11/12	0.98	0.04	23,25,27,28	0
3	BGC	N	2	11/12	0.98	0.04	27,29,30,30	0
2	BGC	K	2	11/12	0.98	0.04	29,30,32,33	0
3	BGC	P	1	12/12	0.98	0.05	25,27,28,32	0
3	BGC	P	2	11/12	0.98	0.04	27,28,29,29	0
2	BGC	Q	2	11/12	0.99	0.03	23,24,25,26	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.

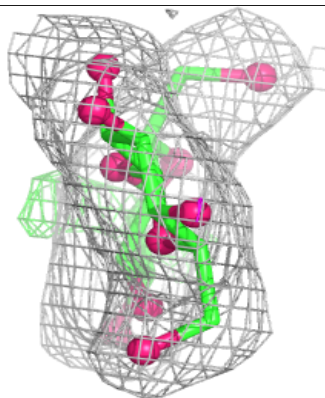
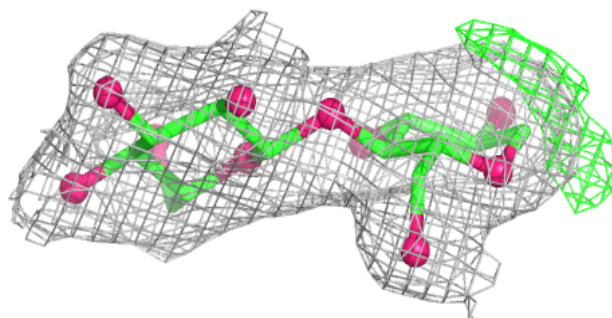
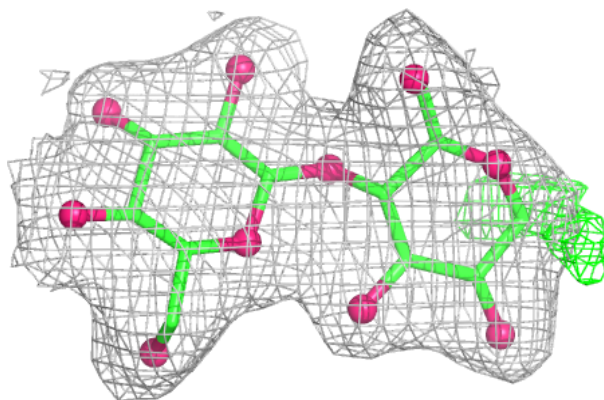


Electron density around Chain I:

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

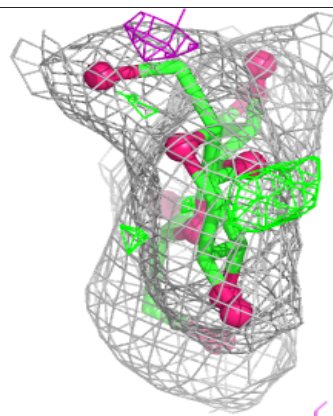
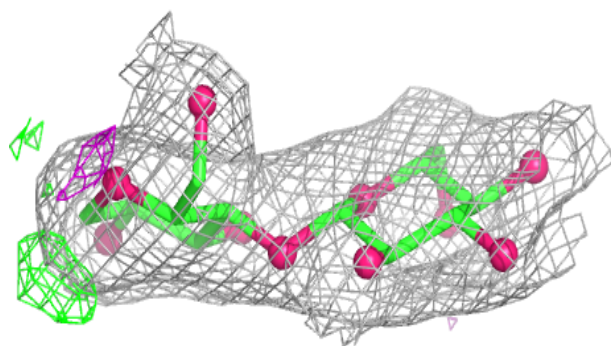
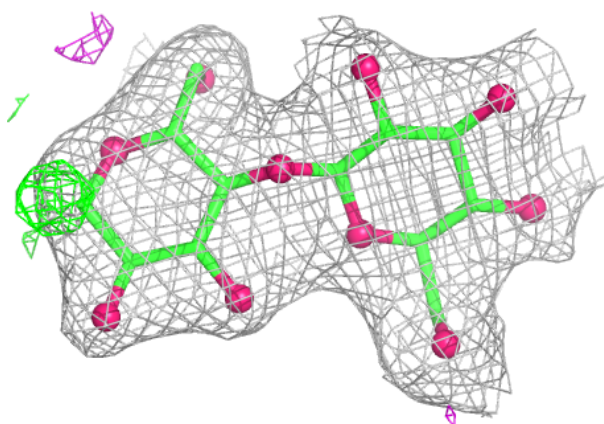
**Electron density around Chain K:**

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



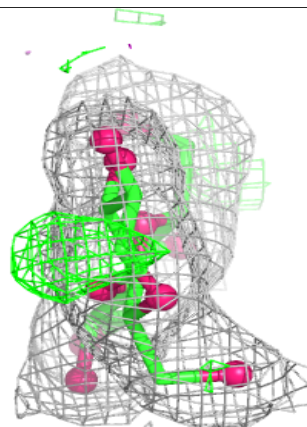
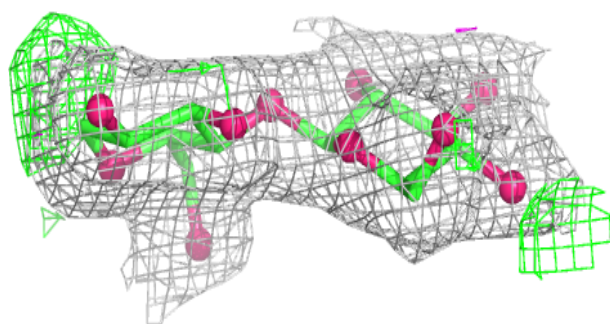
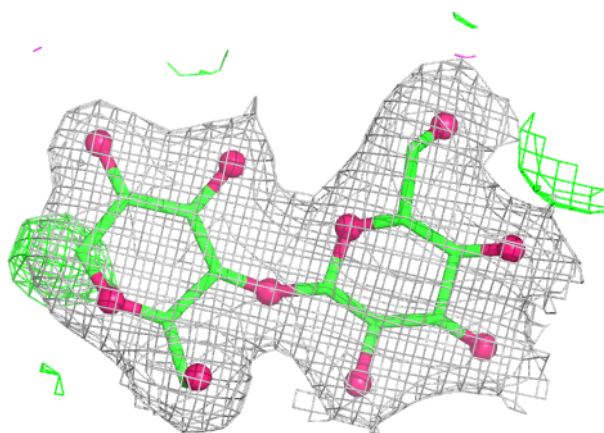
Electron density around Chain M:

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



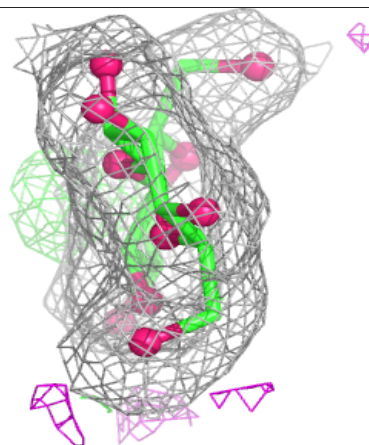
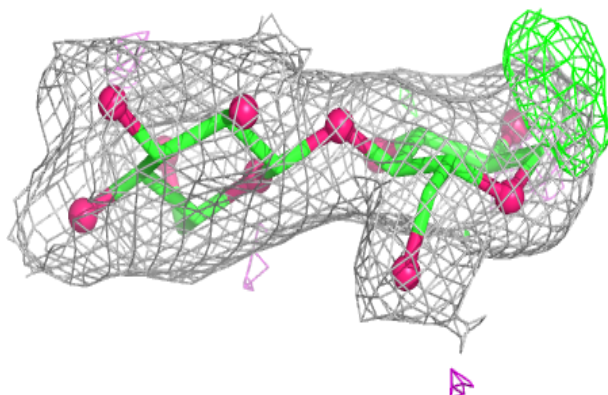
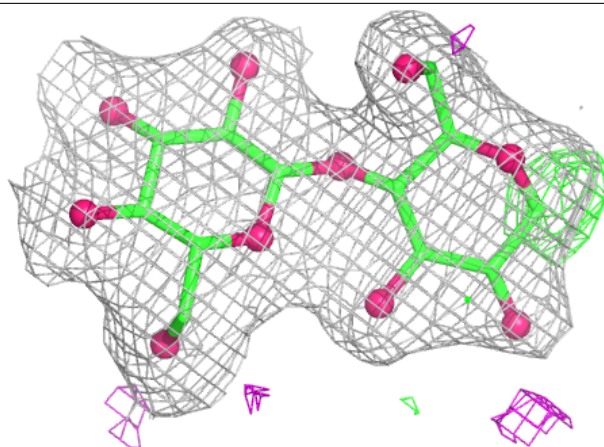
Electron density around Chain O:

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

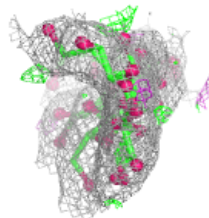
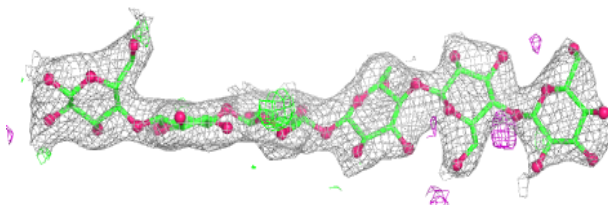
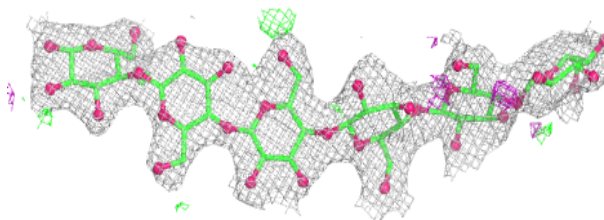


Electron density around Chain Q:

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

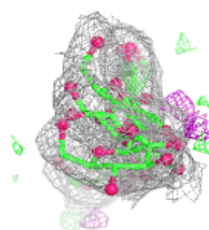
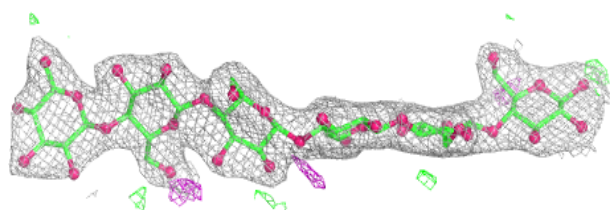
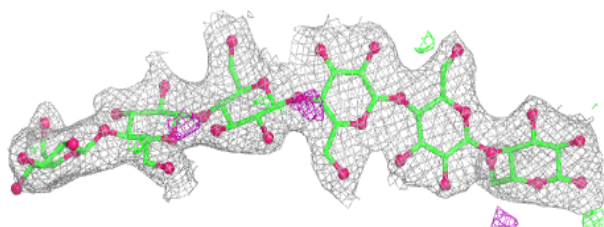
**Electron density around Chain H:**

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

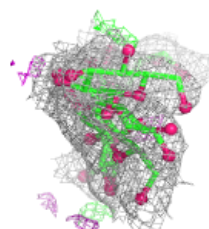
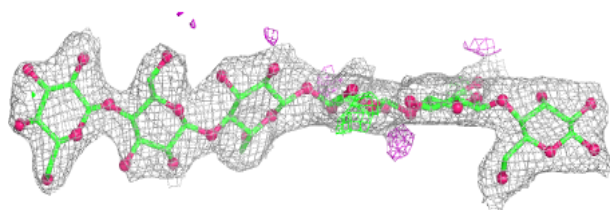
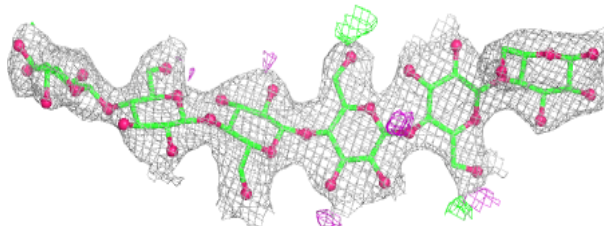


Electron density around Chain J:

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

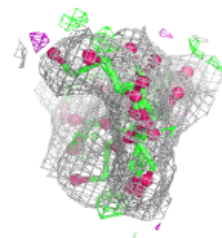
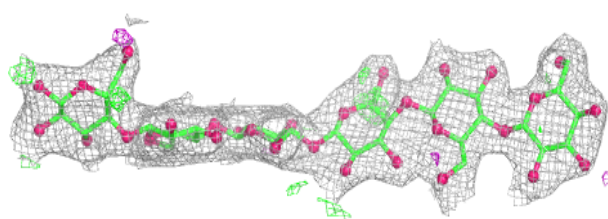
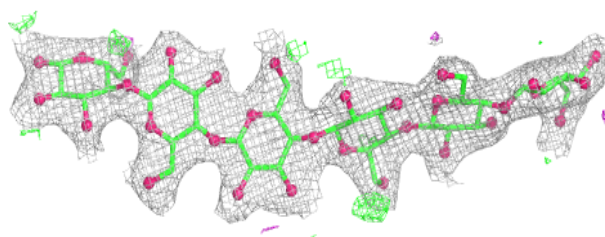
**Electron density around Chain L:**

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

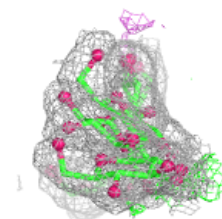
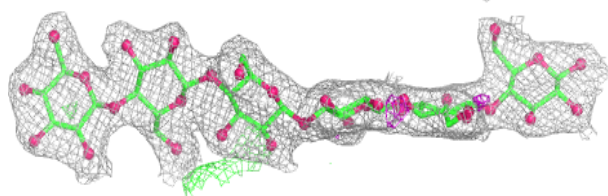
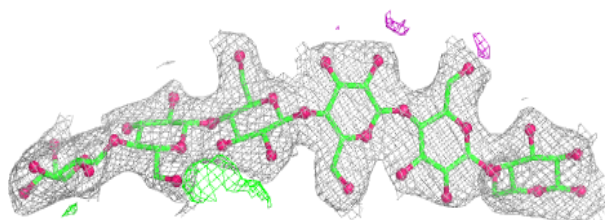


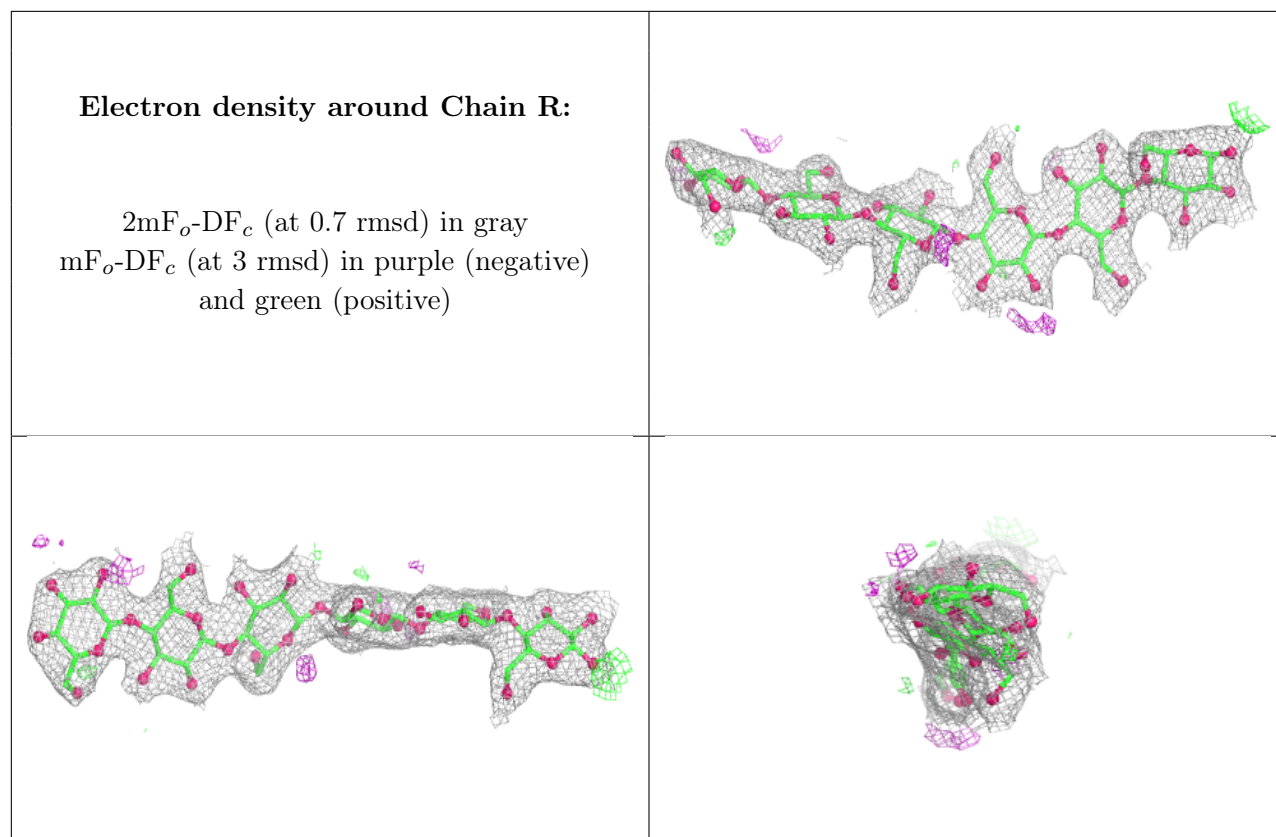
Electron density around Chain N:

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

**Electron density around Chain P:**

$2mF_o-DF_c$ (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.