



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 20, 2026 – 04:39 AM UTC

PDB ID : 1J10 / pdb_00001j10
Title : beta-amylase from *Bacillus cereus* var. *mycoides* in complex with GGX
Authors : Oyama, T.; Miyake, H.; Kusunoki, M.; Nitta, Y.
Deposited on : 2002-11-25
Resolution : 2.10 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

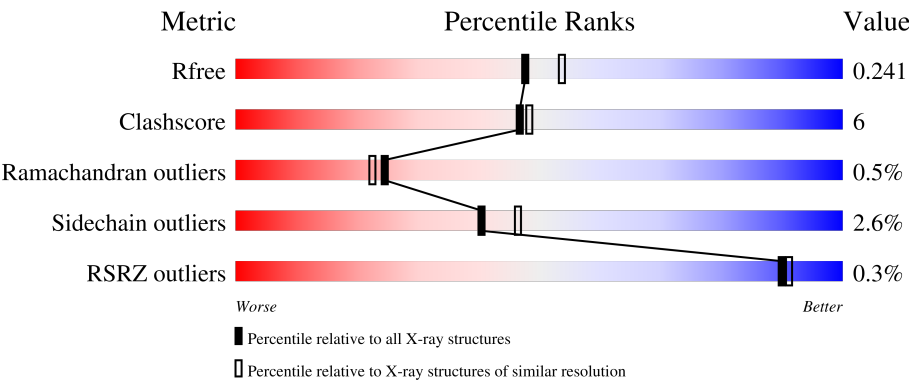
MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.10 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.



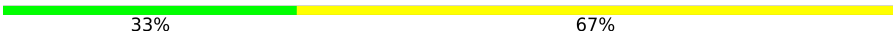
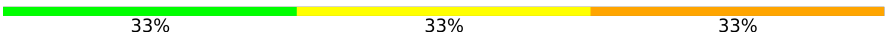
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	6658 (2.10-2.10)
Clashscore	190562	7164 (2.10-2.10)
Ramachandran outliers	187476	7099 (2.10-2.10)
Sidechain outliers	187428	7100 (2.10-2.10)
RSRZ outliers	180081	6662 (2.10-2.10)

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	516	83%16%.
1	B	516	%82%15%.
1	C	516	80%18%.
1	D	516	82%16%.
2	E	3	67%33%

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Mol	Chain	Length	Quality of chain
2	F	3	 67%33%
2	G	3	 100%
2	H	3	 67%33%
2	J	3	 67%33%
2	K	3	 100%
2	L	3	 33%67%
3	I	3	 33%33%33%

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 17298 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 Beta-amylase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	516	Total	C	N	O	S	0	0	0
			4119	2645	676	781	17			
1	B	516	Total	C	N	O	S	0	0	0
			4119	2645	676	781	17			
1	C	516	Total	C	N	O	S	0	0	0
			4119	2645	676	781	17			
1	D	516	Total	C	N	O	S	0	0	0
			4119	2645	676	781	17			

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



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace
2	E	3	Total	C	O	1	0	0
			32	17	15			
2	F	3	Total	C	O	1	0	0
			32	17	15			
2	G	3	Total	C	O	1	0	0
			32	17	15			
2	H	3	Total	C	O	1	0	0
			32	17	15			
2	J	3	Total	C	O	1	0	0
			32	17	15			
2	K	3	Total	C	O	1	0	0
			32	17	15			
2	L	3	Total	C	O	1	0	0
			32	17	15			

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



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace
3	I	3	Total	C	O	1	0	0
			32	17	15			

- Molecule 4 is CALCIUM ION (CCD ID: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	1	Total	Ca	0	0
			1	1		
4	B	1	Total	Ca	0	0
			1	1		
4	C	1	Total	Ca	0	0
			1	1		
4	D	1	Total	Ca	0	0
			1	1		

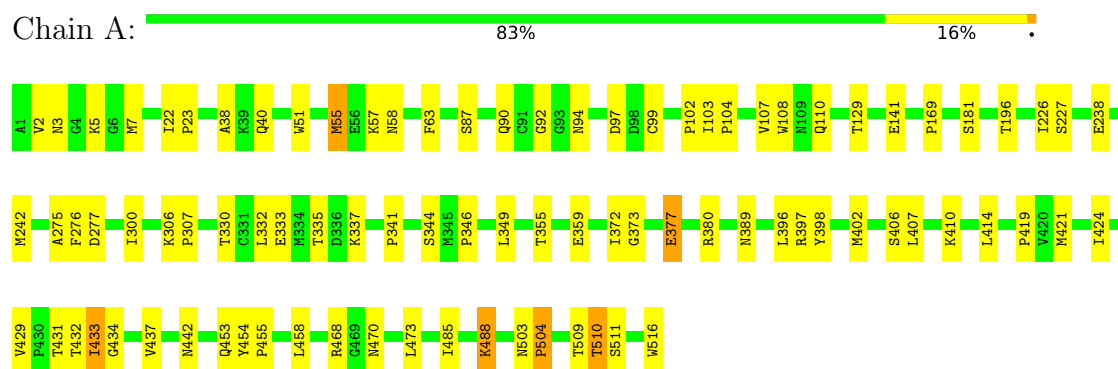
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	160	Total	O	0	0
			160	160		
5	B	120	Total	O	0	0
			120	120		
5	C	128	Total	O	0	0
			128	128		
5	D	154	Total	O	0	0
			154	154		

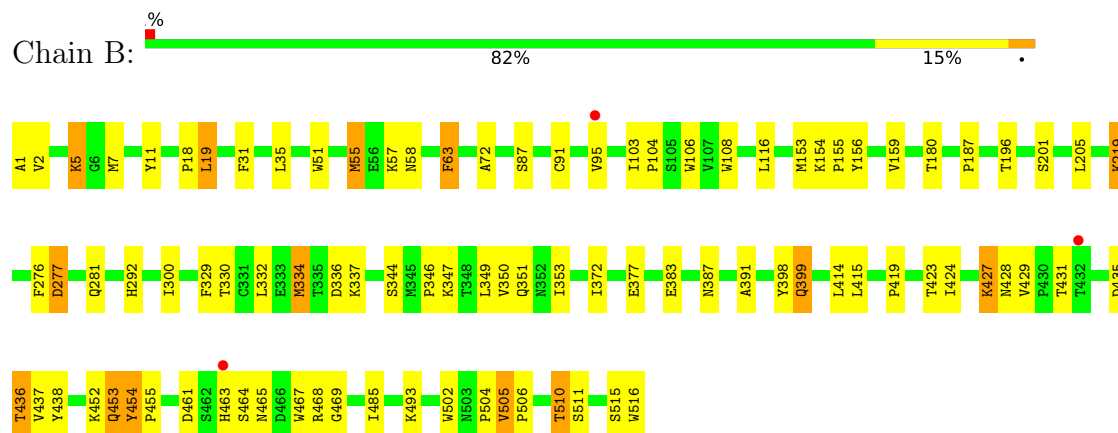
3 Residue-property plots

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.

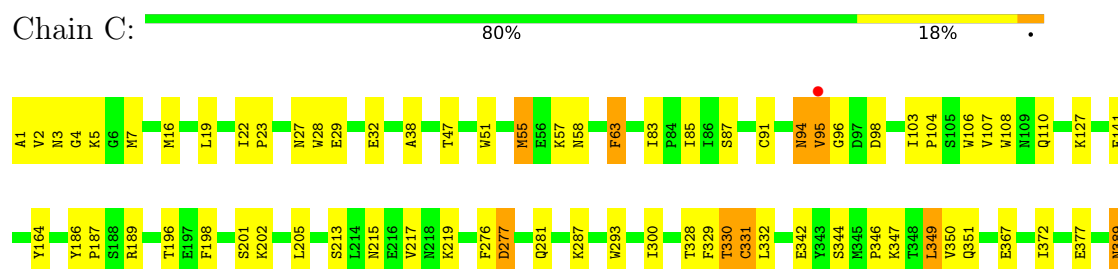
• Molecule 1: Beta-amylase

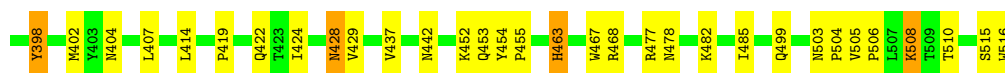


• Molecule 1: Beta-amylase



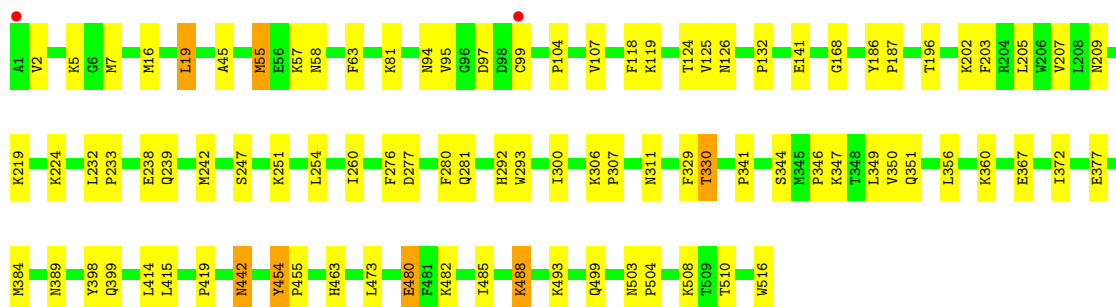
• Molecule 1: Beta-amylase





- Molecule 1: Beta-amylase

Chain D: 82% 16%



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

Chain E: 67% 33%



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

Chain F: 67% 33%



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

Chain G: 100%



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

Chain H: 67% 33%



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

Chain J:  67% 33%



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

Chain K:  100%



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

Chain L:  33% 67%



- Molecule 3: alpha-D-glucopyranose-(1-4)-alpha-D-glucopyranose-(1-4)-alpha-D-xylopyranose

Chain I:  33% 33% 33%



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	177.90Å 112.90Å 146.20Å 90.00° 105.80° 90.00°	Depositor
Resolution (Å)	8.00 – 2.10 8.00 – 2.10	Depositor EDS
% Data completeness (in resolution range)	85.7 (8.00-2.10) 84.0 (8.00-2.10)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.89 (at 2.00Å)	Xtriage
Refinement program	X-PLOR 3.1	Depositor
R, R_{free}	0.189 , 0.239 0.196 , 0.241	Depositor DCC
R_{free} test set	7630 reflections (4.95%)	wwPDB-VP
Wilson B-factor (Å ²)	23.5	Xtriage
Anisotropy	0.216	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 64.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	17298	wwPDB-VP
Average B, all atoms (Å ²)	28.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.38% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: YYS, CA, XYP, GLC

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	0.43	0/4234	0.92	17/5751 (0.3%)
1	B	0.43	0/4234	0.91	17/5751 (0.3%)
1	C	0.44	0/4234	0.91	19/5751 (0.3%)
1	D	0.42	0/4234	0.90	16/5751 (0.3%)
All	All	0.43	0/16936	0.91	69/23004 (0.3%)

There are no bond length outliers.

All (69) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	332	LEU	N-CA-C	8.76	121.98	111.82
1	A	442	ASN	N-CA-C	8.57	120.40	111.14
1	D	330	THR	N-CA-C	8.45	122.39	112.93
1	B	464	SER	N-CA-C	-8.14	103.35	113.28
1	C	58	ASN	N-CA-C	8.04	121.14	111.82
1	D	57	LYS	N-CA-C	8.04	119.67	111.07
1	B	332	LEU	N-CA-C	8.02	122.32	112.54
1	A	277	ASP	N-CA-C	7.49	120.16	111.02
1	B	57	LYS	N-CA-C	7.38	118.97	111.07
1	D	442	ASN	N-CA-C	7.29	119.22	111.28
1	A	337	LYS	N-CA-C	-7.13	102.57	112.12
1	A	57	LYS	N-CA-C	6.80	118.38	110.97
1	C	55	MET	N-CA-C	6.79	119.70	111.82
1	C	442	ASN	N-CA-C	6.75	118.72	111.36
1	B	330	THR	N-CA-C	6.74	120.79	111.55
1	B	505	VAL	N-CA-C	-6.58	102.50	108.95
1	A	276	PHE	N-CA-C	6.49	121.33	113.41
1	C	389	ASN	N-CA-C	6.43	120.43	111.74
1	C	277	ASP	N-CA-C	6.30	118.70	111.02
1	B	58	ASN	N-CA-C	6.22	120.13	112.54

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	196	THR	N-CA-C	-6.21	102.34	110.53
1	D	277	ASP	N-CA-C	6.21	118.56	111.11
1	C	342	GLU	N-CA-C	6.18	118.02	111.28
1	A	414	LEU	N-CA-C	6.16	118.92	111.40
1	B	55	MET	N-CA-C	6.15	120.05	112.54
1	A	332	LEU	N-CA-C	6.04	120.92	113.50
1	C	141	GLU	N-CA-C	6.03	117.52	111.07
1	D	454	TYR	N-CA-C	6.01	123.10	109.81
1	A	58	ASN	N-CA-C	6.01	119.87	112.54
1	C	87	SER	N-CA-C	5.99	119.26	108.17
1	C	57	LYS	N-CA-C	5.96	117.47	110.97
1	A	55	MET	N-CA-C	5.95	120.27	112.89
1	C	330	THR	N-CA-C	5.91	118.37	107.73
1	D	196	THR	N-CA-C	-5.90	101.79	110.52
1	D	292	HIS	N-CA-C	5.89	119.72	112.54
1	D	55	MET	N-CA-C	5.87	118.63	111.82
1	D	389	ASN	N-CA-C	5.87	119.75	112.24
1	A	196	THR	N-CA-C	-5.84	102.67	110.55
1	C	414	LEU	N-CA-C	5.84	118.87	111.69
1	D	58	ASN	N-CA-C	5.72	119.51	112.54
1	A	87	SER	N-CA-C	5.70	118.71	108.17
1	B	277	ASP	N-CA-C	5.69	117.97	111.02
1	C	452	LYS	N-CA-C	5.69	118.58	111.24
1	D	276	PHE	N-CA-C	5.67	120.85	113.88
1	C	276	PHE	N-CA-C	5.63	121.18	113.97
1	A	333	GLU	N-CA-C	5.52	119.86	113.18
1	B	415	LEU	N-CA-C	5.49	120.85	114.04
1	D	141	GLU	N-CA-C	5.43	116.88	111.07
1	A	275	ALA	N-CA-C	5.38	117.22	111.36
1	A	330	THR	N-CA-C	5.36	120.01	112.72
1	B	414	LEU	N-CA-C	5.30	117.86	111.40
1	A	389	ASN	N-CA-C	5.30	119.02	112.24
1	A	51	TRP	N-CA-C	5.29	117.90	109.96
1	C	429	VAL	CA-C-N	5.22	126.37	119.84
1	C	429	VAL	C-N-CA	5.22	126.37	119.84
1	C	189	ARG	N-CA-C	-5.22	103.64	110.53
1	A	141	GLU	N-CA-C	5.20	116.95	111.28
1	D	168	GLY	CA-C-N	5.19	125.10	119.76
1	D	168	GLY	C-N-CA	5.19	125.10	119.76
1	B	116	LEU	N-CA-C	5.18	119.32	112.89
1	B	334	MET	N-CA-C	5.18	118.45	110.42
1	D	414	LEU	N-CA-C	5.17	118.06	111.69

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	196	THR	N-CA-C	-5.10	103.66	110.55
1	B	63	PHE	N-CA-C	5.09	118.01	110.48
1	B	87	SER	N-CA-C	5.05	117.52	108.17
1	B	276	PHE	N-CA-C	5.02	120.39	113.97
1	C	63	PHE	N-CA-C	5.01	117.90	110.48
1	B	292	HIS	N-CA-C	5.01	119.39	113.28
1	D	415	LEU	N-CA-C	5.00	120.76	114.31

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	4119	0	3984	42	0
1	B	4119	0	3984	49	0
1	C	4119	0	3984	51	0
1	D	4119	0	3984	43	0
2	E	32	0	19	0	0
2	F	32	0	19	0	0
2	G	32	0	19	0	0
2	H	32	0	19	0	0
2	J	32	0	19	0	0
2	K	32	0	19	0	0
2	L	32	0	19	0	0
3	I	32	0	27	2	0
4	A	1	0	0	0	0
4	B	1	0	0	0	0
4	C	1	0	0	0	0
4	D	1	0	0	0	0
5	A	160	0	0	1	0
5	B	120	0	0	1	0
5	C	128	0	0	0	0
5	D	154	0	0	0	0
All	All	17298	0	16096	187	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 (187) 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)
3:I:1:XY5:H52	3:I:2:GLC:H5	1.50	0.92
1:A:7:MET:HE2	1:A:419:PRO:HB3	1.63	0.80
1:A:238:GLU:HG2	1:A:242:MET:HE2	1.67	0.77
1:B:277:ASP:O	1:B:281:GLN:HG3	1.90	0.71
3:I:1:XY5:H52	3:I:2:GLC:C5	2.21	0.70
1:B:329:PHE:CE2	1:B:350:VAL:HG22	2.27	0.70
1:D:280:PHE:O	1:D:281:GLN:HG2	1.92	0.69
1:D:7:MET:HE2	1:D:419:PRO:HG3	1.74	0.68
1:D:2:VAL:O	1:D:5:LYS:HG2	1.94	0.67
1:A:103:ILE:HD11	1:A:108:TRP:CZ2	2.33	0.64
1:B:2:VAL:O	1:B:5:LYS:HG3	1.99	0.63
1:D:247:SER:O	1:D:251:LYS:HG3	1.99	0.62
1:A:421:MET:SD	1:A:470:ASN:HB3	2.41	0.60
1:A:488:LYS:HD3	1:A:488:LYS:H	1.66	0.60
1:B:431:THR:OG1	1:B:435:ASP:HB2	2.02	0.60
1:B:1:ALA:HA	1:B:387:ASN:OD1	2.00	0.59
1:C:372:ILE:HG23	1:C:377:GLU:HB2	1.81	0.59
1:A:434:GLY:HA3	1:A:488:LYS:HG3	1.85	0.59
1:B:201:SER:O	1:B:205:LEU:HG	2.01	0.59
1:C:198:PHE:O	1:C:202:LYS:HG2	2.02	0.59
1:A:355:THR:O	1:A:359:GLU:HG3	2.02	0.58
1:B:329:PHE:CE1	1:B:353:ILE:HD12	2.39	0.58
1:D:205:LEU:O	1:D:209:ASN:HB2	2.04	0.58
1:D:346:PRO:O	1:D:350:VAL:HG23	2.04	0.58
1:B:7:MET:HE2	1:B:419:PRO:HB3	1.85	0.58
1:D:351:GLN:HG2	1:D:384:MET:HE1	1.85	0.58
1:C:4:GLY:HA2	1:C:455:PRO:O	2.04	0.57
1:A:2:VAL:O	1:A:5:LYS:HG2	2.03	0.57
1:C:201:SER:O	1:C:205:LEU:HG	2.04	0.57
1:C:422:GLN:NE2	1:C:506:PRO:HG2	2.20	0.57
1:D:372:ILE:HG23	1:D:377:GLU:HB2	1.87	0.56
1:A:488:LYS:HD3	1:A:488:LYS:N	2.21	0.56
1:B:156:TYR:O	1:B:159:VAL:HG22	2.06	0.55
1:D:119:LYS:HA	1:D:124:THR:O	2.05	0.55
1:B:424:ILE:O	1:B:468:ARG:HA	2.06	0.55
1:D:488:LYS:HA	1:D:488:LYS:HE3	1.88	0.55
1:C:454:TYR:N	1:C:455:PRO:HD2	2.22	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:103:ILE:HD11	1:B:108:TRP:CZ2	2.42	0.54
1:D:238:GLU:HG2	1:D:242:MET:HE2	1.90	0.54
1:B:337:LYS:HD2	1:B:344:SER:HB3	1.89	0.53
1:B:372:ILE:HG23	1:B:377:GLU:HB2	1.90	0.53
1:C:1:ALA:HB3	1:C:5:LYS:HG3	1.90	0.53
1:B:300:ILE:HG13	1:B:300:ILE:O	2.08	0.53
1:A:226:ILE:HD12	1:A:227:SER:N	2.23	0.52
1:A:488:LYS:H	1:A:488:LYS:CD	2.22	0.52
1:A:107:VAL:O	1:A:110:GLN:HG2	2.09	0.52
1:D:104:PRO:HB2	1:D:107:VAL:HG23	1.92	0.52
1:A:300:ILE:O	1:A:300:ILE:HG13	2.09	0.52
1:C:300:ILE:HG13	1:C:300:ILE:O	2.09	0.52
1:D:329:PHE:CE2	1:D:350:VAL:HG22	2.44	0.52
1:C:27:ASN:OD1	1:C:29:GLU:HB3	2.09	0.52
1:C:28:TRP:O	1:C:32:GLU:HG3	2.10	0.52
1:D:356:LEU:HB3	1:D:360:LYS:NZ	2.25	0.52
1:C:477:ARG:O	1:C:505:VAL:HG23	2.09	0.51
1:B:461:ASP:O	1:B:465:ASN:N	2.43	0.51
1:C:5:LYS:HE3	1:C:389:ASN:HB2	1.93	0.51
1:B:219:LYS:HB3	1:B:219:LYS:HZ3	1.75	0.51
1:B:427:LYS:HB3	1:B:427:LYS:NZ	2.26	0.51
1:B:11:TYR:HA	1:B:391:ALA:O	2.11	0.51
1:C:346:PRO:O	1:C:350:VAL:HG23	2.11	0.51
1:C:347:LYS:O	1:C:351:GLN:HG3	2.10	0.50
1:C:2:VAL:O	1:C:5:LYS:HG2	2.12	0.50
1:A:55:MET:O	1:A:63:PHE:HA	2.13	0.49
1:B:1:ALA:HB1	1:B:5:LYS:HE2	1.93	0.49
1:A:454:TYR:N	1:A:455:PRO:HD2	2.27	0.49
1:C:47:THR:HG22	1:C:83:ILE:HB	1.93	0.49
1:C:453:GLN:C	1:C:455:PRO:HD2	2.38	0.49
1:C:330:THR:HG22	1:C:367:GLU:CG	2.43	0.49
1:D:118:PHE:HB2	1:D:126:ASN:HB3	1.93	0.49
1:A:306:LYS:HB2	1:A:307:PRO:HD3	1.94	0.49
1:A:424:ILE:O	1:A:468:ARG:HA	2.12	0.49
1:C:404:ASN:HD21	1:C:407:LEU:HD23	1.77	0.49
1:B:454:TYR:N	1:B:455:PRO:HD2	2.28	0.48
1:C:103:ILE:HD11	1:C:108:TRP:CZ2	2.48	0.48
1:B:383:GLU:O	1:B:387:ASN:HB2	2.13	0.48
1:A:485:ILE:HG13	1:A:516:TRP:CH2	2.48	0.48
1:D:97:ASP:C	1:D:99:CYS:H	2.22	0.48
1:D:485:ILE:HG13	1:D:516:TRP:CH2	2.48	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:437:VAL:HB	1:B:467:TRP:CE3	2.49	0.48
1:A:433:ILE:HD12	1:A:433:ILE:H	1.79	0.48
1:B:154:LYS:HB3	1:B:155:PRO:HD3	1.96	0.48
1:B:429:VAL:HG22	1:B:516:TRP:CE3	2.49	0.48
1:C:7:MET:HE2	1:C:419:PRO:HB3	1.96	0.47
1:A:344:SER:C	1:A:346:PRO:HD3	2.39	0.47
1:C:107:VAL:O	1:C:110:GLN:HG2	2.14	0.47
1:D:300:ILE:HG13	1:D:300:ILE:O	2.13	0.47
1:D:254:LEU:HD13	1:D:311:ASN:ND2	2.29	0.47
1:D:508:LYS:N	1:D:508:LYS:HD2	2.29	0.47
1:C:108:TRP:CZ2	1:C:127:LYS:HB3	2.49	0.47
1:A:429:VAL:HG11	1:A:437:VAL:HG21	1.96	0.47
1:B:19:LEU:HD12	1:B:19:LEU:HA	1.78	0.47
1:C:55:MET:O	1:C:63:PHE:HA	2.15	0.47
1:C:508:LYS:NZ	1:C:508:LYS:HB3	2.29	0.47
1:D:329:PHE:HE2	1:D:350:VAL:HG22	1.79	0.47
1:C:329:PHE:HZ	1:C:349:LEU:HD13	1.78	0.46
1:C:424:ILE:O	1:C:468:ARG:HA	2.14	0.46
1:D:132:PRO:HB2	1:D:260:ILE:HG22	1.96	0.46
1:A:372:ILE:HG23	1:A:377:GLU:HB2	1.97	0.46
1:C:277:ASP:O	1:C:281:GLN:HG2	2.15	0.46
1:B:454:TYR:N	1:B:455:PRO:CD	2.79	0.46
1:C:38:ALA:HB2	1:C:402:MET:CE	2.46	0.46
1:C:85:ILE:HG12	1:C:164:TYR:HB2	1.98	0.46
1:D:239:GLN:HA	1:D:242:MET:HE3	1.97	0.46
1:D:55:MET:O	1:D:63:PHE:HA	2.15	0.46
1:C:404:ASN:ND2	1:C:407:LEU:HD23	2.31	0.46
1:B:428:ASN:HD22	1:B:515:SER:HB3	1.81	0.46
1:C:104:PRO:HB3	1:C:106:TRP:NE1	2.31	0.46
1:A:432:THR:HG22	1:A:433:ILE:HD12	1.98	0.46
1:B:51:TRP:CZ2	1:B:91:CYS:SG	3.09	0.46
1:B:334:MET:HE2	1:B:337:LYS:NZ	2.31	0.46
1:D:238:GLU:O	1:D:242:MET:HG3	2.16	0.46
1:B:485:ILE:HG13	1:B:516:TRP:CH2	2.50	0.45
1:A:94:ASN:HB2	1:A:97:ASP:OD1	2.16	0.45
1:B:180:THR:HG22	5:B:791:HOH:O	2.16	0.45
1:B:219:LYS:HB3	1:B:219:LYS:NZ	2.31	0.45
1:B:502:TRP:CE2	1:B:506:PRO:HD3	2.51	0.45
1:D:454:TYR:N	1:D:455:PRO:HD2	2.31	0.45
1:C:454:TYR:N	1:C:455:PRO:CD	2.79	0.45
1:D:344:SER:C	1:D:346:PRO:HD3	2.41	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:92:GLY:HA3	1:A:99:CYS:O	2.17	0.45
1:D:347:LYS:O	1:D:351:GLN:HG3	2.17	0.45
1:B:72:ALA:HB2	1:B:153:MET:HE1	1.99	0.45
1:C:344:SER:C	1:C:346:PRO:HD3	2.42	0.45
1:A:129:THR:HG21	1:A:169:PRO:HD3	1.98	0.45
1:A:373:GLY:HA2	1:A:407:LEU:HD21	1.98	0.45
1:B:55:MET:O	1:B:63:PHE:HA	2.18	0.44
1:C:215:ASN:O	1:C:219:LYS:HG3	2.17	0.44
1:A:437:VAL:HG12	1:A:458:LEU:HD12	1.99	0.44
1:A:104:PRO:HB2	1:A:107:VAL:HG23	2.00	0.44
1:C:213:SER:O	1:C:217:VAL:HG23	2.18	0.44
1:C:478:ASN:HB3	1:C:503:ASN:HD22	1.83	0.44
1:B:344:SER:C	1:B:346:PRO:HD3	2.43	0.44
1:B:399:GLN:H	1:B:399:GLN:CD	2.25	0.44
1:C:437:VAL:HB	1:C:467:TRP:CE3	2.53	0.44
1:B:423:THR:HA	1:B:469:GLY:O	2.18	0.44
1:C:186:TYR:HD1	1:C:293:TRP:CD2	2.36	0.44
1:D:330:THR:HA	1:D:367:GLU:HB3	1.99	0.44
1:A:396:LEU:HA	1:A:397:ARG:HA	1.80	0.43
1:A:406:SER:O	1:A:410:LYS:HG2	2.18	0.43
1:A:454:TYR:N	1:A:455:PRO:CD	2.81	0.43
1:C:19:LEU:HD23	1:C:98:ASP:HB3	1.99	0.43
1:A:22:ILE:N	1:A:23:PRO:HD2	2.33	0.43
1:A:432:THR:HG22	1:A:433:ILE:N	2.33	0.43
1:D:19:LEU:HD12	1:D:19:LEU:HA	1.82	0.43
1:B:510:THR:HB	1:B:511:SER:H	1.53	0.43
1:B:104:PRO:HB3	1:B:106:TRP:NE1	2.34	0.43
1:C:51:TRP:CZ2	1:C:91:CYS:SG	3.12	0.43
1:C:428:ASN:HD21	1:C:515:SER:HB3	1.84	0.43
1:D:203:PHE:O	1:D:207:VAL:HG23	2.19	0.43
1:A:335:THR:HG23	1:A:380:ARG:HD3	2.01	0.42
1:A:510:THR:HB	1:A:511:SER:H	1.69	0.42
1:C:463:HIS:ND1	1:C:463:HIS:C	2.77	0.42
1:B:31:PHE:O	1:B:35:LEU:HG	2.19	0.42
1:A:421:MET:O	1:A:509:THR:HG23	2.20	0.42
1:A:503:ASN:HA	1:A:504:PRO:HA	1.78	0.42
1:A:2:VAL:HG12	1:A:3:ASN:ND2	2.34	0.42
1:D:224:LYS:HA	1:D:224:LYS:HD2	1.72	0.42
1:C:16:MET:O	1:C:398:TYR:HB2	2.20	0.42
1:C:94:ASN:HB3	1:C:95:VAL:H	1.65	0.42
1:B:347:LYS:O	1:B:351:GLN:HG3	2.20	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:436:THR:HB	1:B:438:TYR:CE1	2.55	0.42
1:C:22:ILE:HB	1:C:23:PRO:HD3	2.01	0.42
1:D:485:ILE:HG21	1:D:493:LYS:HD3	2.02	0.41
1:C:2:VAL:O	1:C:3:ASN:HB2	2.20	0.41
1:A:453:GLN:HA	5:A:809:HOH:O	2.21	0.41
1:B:18:PRO:HB3	1:B:399:GLN:HE22	1.86	0.41
1:D:45:ALA:HB2	1:D:81:LYS:HB2	2.03	0.41
1:D:306:LYS:HB2	1:D:307:PRO:HD3	2.02	0.41
1:A:38:ALA:HB2	1:A:402:MET:HE1	2.03	0.41
1:D:202:LYS:HD3	1:D:202:LYS:HA	1.88	0.41
1:C:287:LYS:HG2	1:C:328:THR:HB	2.03	0.41
1:B:2:VAL:HA	1:B:453:GLN:O	2.21	0.41
1:B:336:ASP:HA	1:B:346:PRO:HD2	2.02	0.41
1:D:186:TYR:HD1	1:D:293:TRP:CD2	2.38	0.41
1:D:232:LEU:HB3	1:D:233:PRO:HD2	2.03	0.41
1:D:442:ASN:ND2	1:D:480:GLU:HG3	2.36	0.41
1:B:219:LYS:HD2	1:B:219:LYS:C	2.46	0.41
1:C:482:LYS:HG3	1:C:499:GLN:HA	2.01	0.41
1:C:485:ILE:HG13	1:C:516:TRP:CH2	2.56	0.41
1:B:493:LYS:HB3	1:B:493:LYS:HE2	1.88	0.40
1:B:1:ALA:CB	1:B:5:LYS:HE2	2.51	0.40
1:D:94:ASN:HB2	1:D:97:ASP:OD2	2.20	0.40
1:D:482:LYS:HG3	1:D:499:GLN:HA	2.02	0.40
1:C:94:ASN:O	1:C:96:GLY:N	2.54	0.40
1:D:118:PHE:O	1:D:125:VAL:HA	2.22	0.40
1:D:503:ASN:HA	1:D:504:PRO:HA	1.81	0.40
1:A:90:GLN:HG3	1:A:102:PRO:HA	2.03	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	514/516 (100%)	489 (95%)	24 (5%)	1 (0%)	43	44
1	B	514/516 (100%)	486 (95%)	24 (5%)	4 (1%)	16	12
1	C	514/516 (100%)	486 (95%)	25 (5%)	3 (1%)	21	18
1	D	514/516 (100%)	488 (95%)	24 (5%)	2 (0%)	30	28
All	All	2056/2064 (100%)	1949 (95%)	97 (5%)	10 (0%)	24	22

All (10) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	398	TYR
1	C	95	VAL
1	C	331	CYS
1	B	463	HIS
1	D	95	VAL
1	C	398	TYR
1	D	398	TYR
1	A	398	TYR
1	B	95	VAL
1	B	454	TYR

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	440/440 (100%)	429 (98%)	11 (2%)	42	48
1	B	440/440 (100%)	427 (97%)	13 (3%)	36	41
1	C	440/440 (100%)	431 (98%)	9 (2%)	48	56
1	D	440/440 (100%)	428 (97%)	12 (3%)	39	45
All	All	1760/1760 (100%)	1715 (97%)	45 (3%)	40	46

All (45) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	40	GLN

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Mol	Chain	Res	Type
1	A	181	SER
1	A	341	PRO
1	A	349	LEU
1	A	377	GLU
1	A	431	THR
1	A	433	ILE
1	A	473	LEU
1	A	488	LYS
1	A	504	PRO
1	A	510	THR
1	B	5	LYS
1	B	19	LEU
1	B	187	PRO
1	B	219	LYS
1	B	349	LEU
1	B	399	GLN
1	B	427	LYS
1	B	436	THR
1	B	452	LYS
1	B	453	GLN
1	B	504	PRO
1	B	505	VAL
1	B	510	THR
1	C	94	ASN
1	C	187	PRO
1	C	331	CYS
1	C	349	LEU
1	C	428	ASN
1	C	463	HIS
1	C	504	PRO
1	C	508	LYS
1	C	510	THR
1	D	16	MET
1	D	19	LEU
1	D	187	PRO
1	D	219	LYS
1	D	341	PRO
1	D	349	LEU
1	D	399	GLN
1	D	463	HIS
1	D	473	LEU
1	D	480	GLU

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Mol	Chain	Res	Type
1	D	488	LYS
1	D	510	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (31) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	3	ASN
1	A	40	GLN
1	A	311	ASN
1	A	351	GLN
1	A	352	ASN
1	A	453	GLN
1	A	478	ASN
1	A	499	GLN
1	A	503	ASN
1	B	3	ASN
1	B	73	GLN
1	B	77	ASN
1	B	100	ASN
1	B	239	GLN
1	B	281	GLN
1	B	311	ASN
1	B	428	ASN
1	C	109	ASN
1	C	239	GLN
1	C	281	GLN
1	C	351	GLN
1	C	352	ASN
1	C	428	ASN
1	C	503	ASN
1	D	62	GLN
1	D	69	GLN
1	D	311	ASN
1	D	351	GLN
1	D	470	ASN
1	D	478	ASN
1	D	503	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 ⓘ

24 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	XYP	E	1	2	10,10,10	0.46	0	14,14,14	0.39	0
2	GLC	E	2	2	11,11,12	0.48	0	15,15,17	0.57	0
2	GLC	E	3	2	11,11,12	0.43	0	15,15,17	0.63	1 (6%)
2	XYP	F	1	2	10,10,10	0.73	0	14,14,14	0.59	0
2	GLC	F	2	2	11,11,12	0.51	0	15,15,17	0.63	0
2	GLC	F	3	2	11,11,12	0.30	0	15,15,17	0.72	1 (6%)
2	XYP	G	1	2	10,10,10	0.61	0	14,14,14	0.39	0
2	GLC	G	2	2	11,11,12	0.45	0	15,15,17	0.50	0
2	GLC	G	3	2	11,11,12	0.48	0	15,15,17	0.55	0
2	XYP	H	1	2	10,10,10	0.68	0	14,14,14	0.46	0
2	GLC	H	2	2	11,11,12	0.53	0	15,15,17	0.65	0
2	GLC	H	3	2	11,11,12	0.36	0	15,15,17	0.67	1 (6%)
3	XYS	I	1	3	10,10,10	1.23	1 (10%)	14,14,14	1.46	2 (14%)
3	GLC	I	2	3	11,11,12	0.55	0	15,15,17	0.57	0
3	GLC	I	3	3	11,11,12	0.43	0	15,15,17	0.67	0
2	XYP	J	1	2	10,10,10	0.64	0	14,14,14	0.42	0
2	GLC	J	2	2	11,11,12	0.50	0	15,15,17	0.61	0
2	GLC	J	3	2	11,11,12	0.34	0	15,15,17	0.69	1 (6%)
2	XYP	K	1	2	10,10,10	0.64	0	14,14,14	0.45	0
2	GLC	K	2	2	11,11,12	0.45	0	15,15,17	0.58	0
2	GLC	K	3	2	11,11,12	0.43	0	15,15,17	0.59	0
2	XYP	L	1	2	10,10,10	0.71	0	14,14,14	0.32	0
2	GLC	L	2	2	11,11,12	0.58	0	15,15,17	0.76	1 (6%)
2	GLC	L	3	2	11,11,12	0.33	0	15,15,17	0.65	1 (6%)

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	XYP	E	1	2	-	-	0/1/1/1
2	GLC	E	2	2	-	0/2/19/22	0/1/1/1
2	GLC	E	3	2	-	0/2/19/22	0/1/1/1
2	XYP	F	1	2	-	-	0/1/1/1
2	GLC	F	2	2	-	1/2/19/22	0/1/1/1
2	GLC	F	3	2	-	0/2/19/22	0/1/1/1
2	XYP	G	1	2	-	-	0/1/1/1
2	GLC	G	2	2	-	0/2/19/22	0/1/1/1
2	GLC	G	3	2	-	1/2/19/22	0/1/1/1
2	XYP	H	1	2	-	-	0/1/1/1
2	GLC	H	2	2	-	0/2/19/22	0/1/1/1
2	GLC	H	3	2	-	0/2/19/22	0/1/1/1
3	XYS	I	1	3	-	-	0/1/1/1
3	GLC	I	2	3	-	0/2/19/22	0/1/1/1
3	GLC	I	3	3	-	0/2/19/22	0/1/1/1
2	XYP	J	1	2	-	-	0/1/1/1
2	GLC	J	2	2	-	0/2/19/22	0/1/1/1
2	GLC	J	3	2	-	0/2/19/22	0/1/1/1
2	XYP	K	1	2	-	-	0/1/1/1
2	GLC	K	2	2	-	0/2/19/22	0/1/1/1
2	GLC	K	3	2	-	0/2/19/22	0/1/1/1
2	XYP	L	1	2	-	-	0/1/1/1
2	GLC	L	2	2	-	1/2/19/22	0/1/1/1
2	GLC	L	3	2	-	0/2/19/22	0/1/1/1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	I	1	XYS	O1-C1	-3.33	1.29	1.39

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	I	1	XYS	O1-C1-O5	4.21	120.70	109.91
3	I	1	XYS	O1-C1-C2	3.04	117.82	108.98
2	F	3	GLC	C1-O5-C5	2.38	115.37	112.19
2	J	3	GLC	C1-O5-C5	2.34	115.32	112.19
2	L	2	GLC	C1-O5-C5	2.32	115.30	112.19

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Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
2	H	3	GLC	C1-O5-C5	2.18	115.11	112.19
2	L	3	GLC	C1-O5-C5	2.07	114.96	112.19
2	E	3	GLC	C1-O5-C5	2.06	114.95	112.19

There are no chirality outliers.

All (3) torsion outliers are listed below:

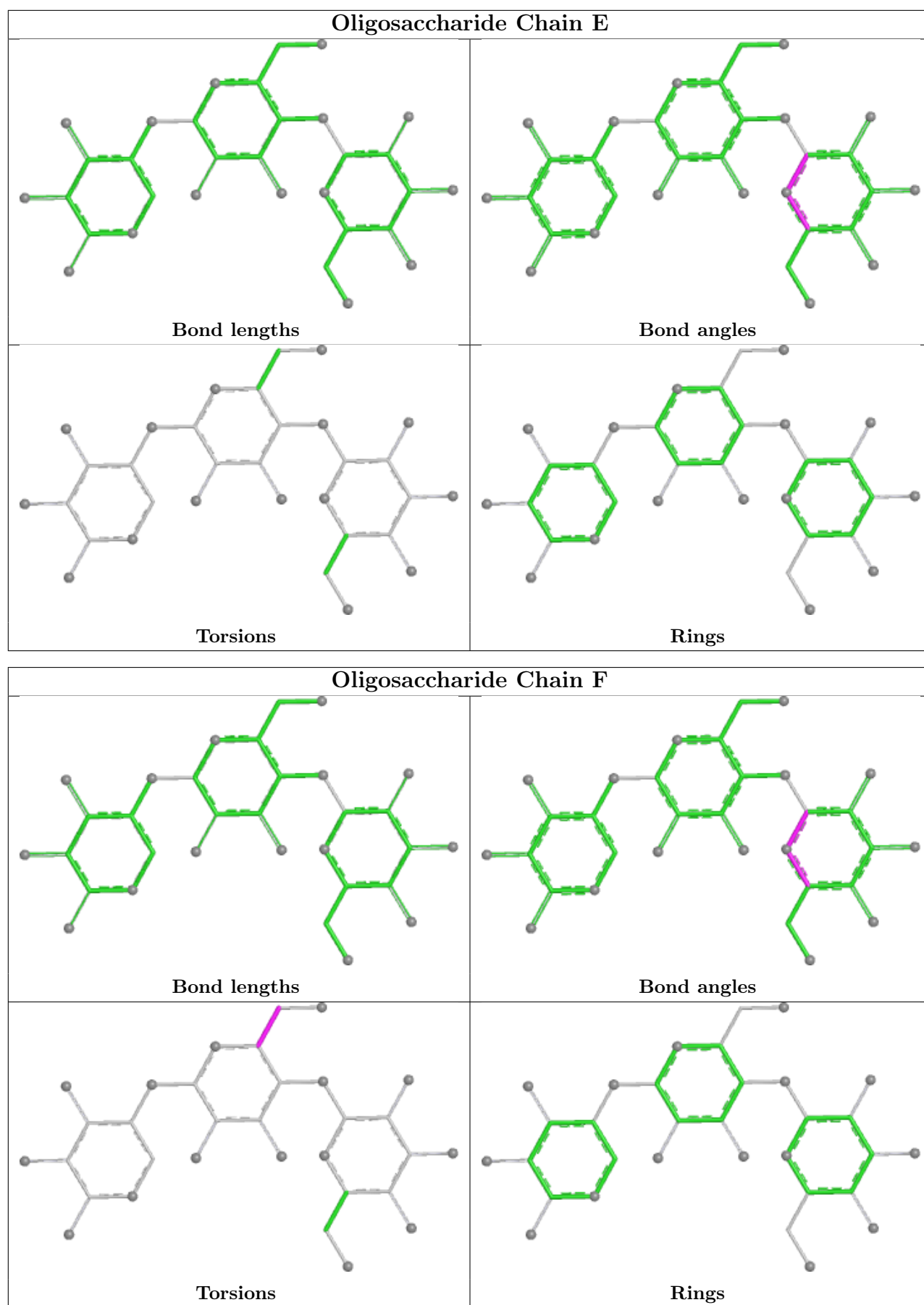
Mol	Chain	Res	Type	Atoms
2	F	2	GLC	O5-C5-C6-O6
2	L	2	GLC	O5-C5-C6-O6
2	G	3	GLC	C4-C5-C6-O6

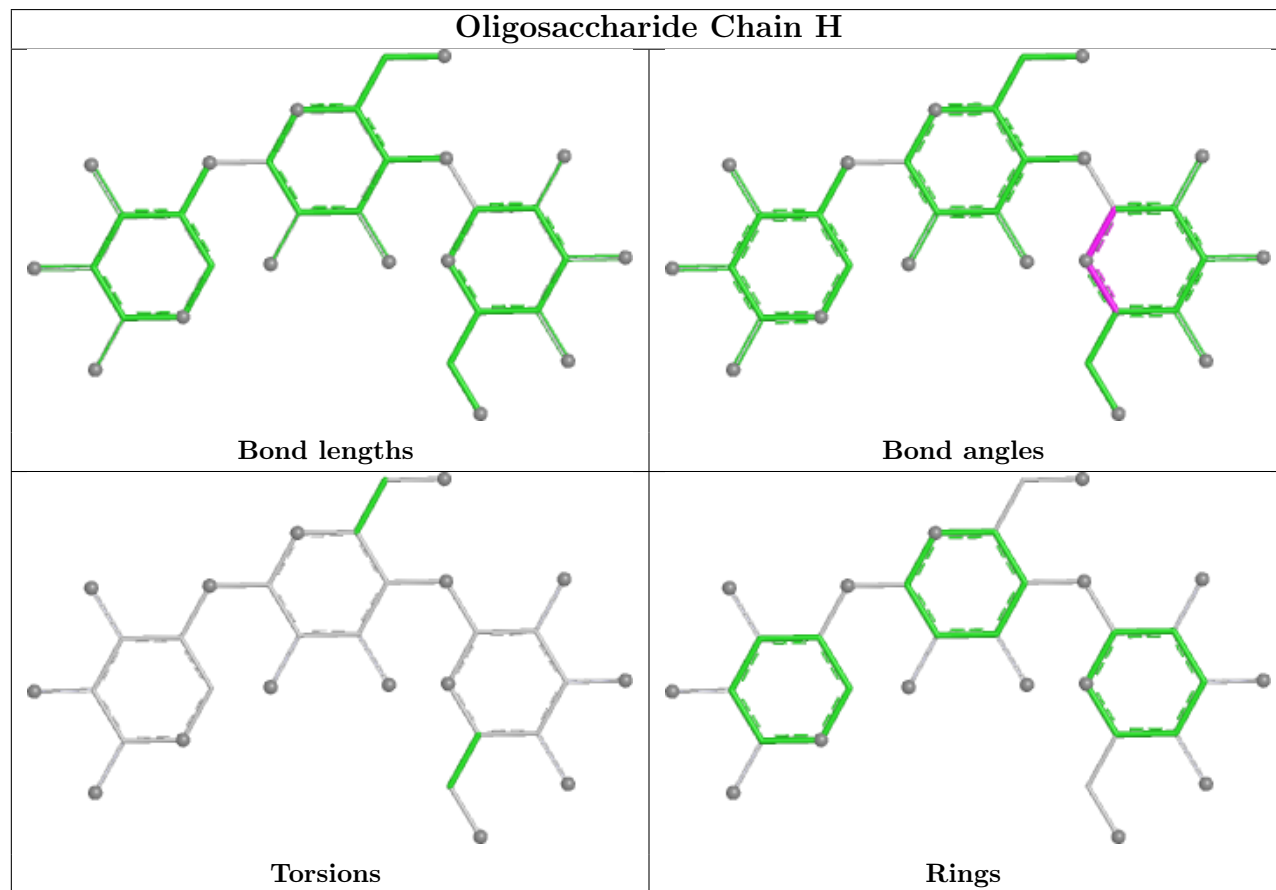
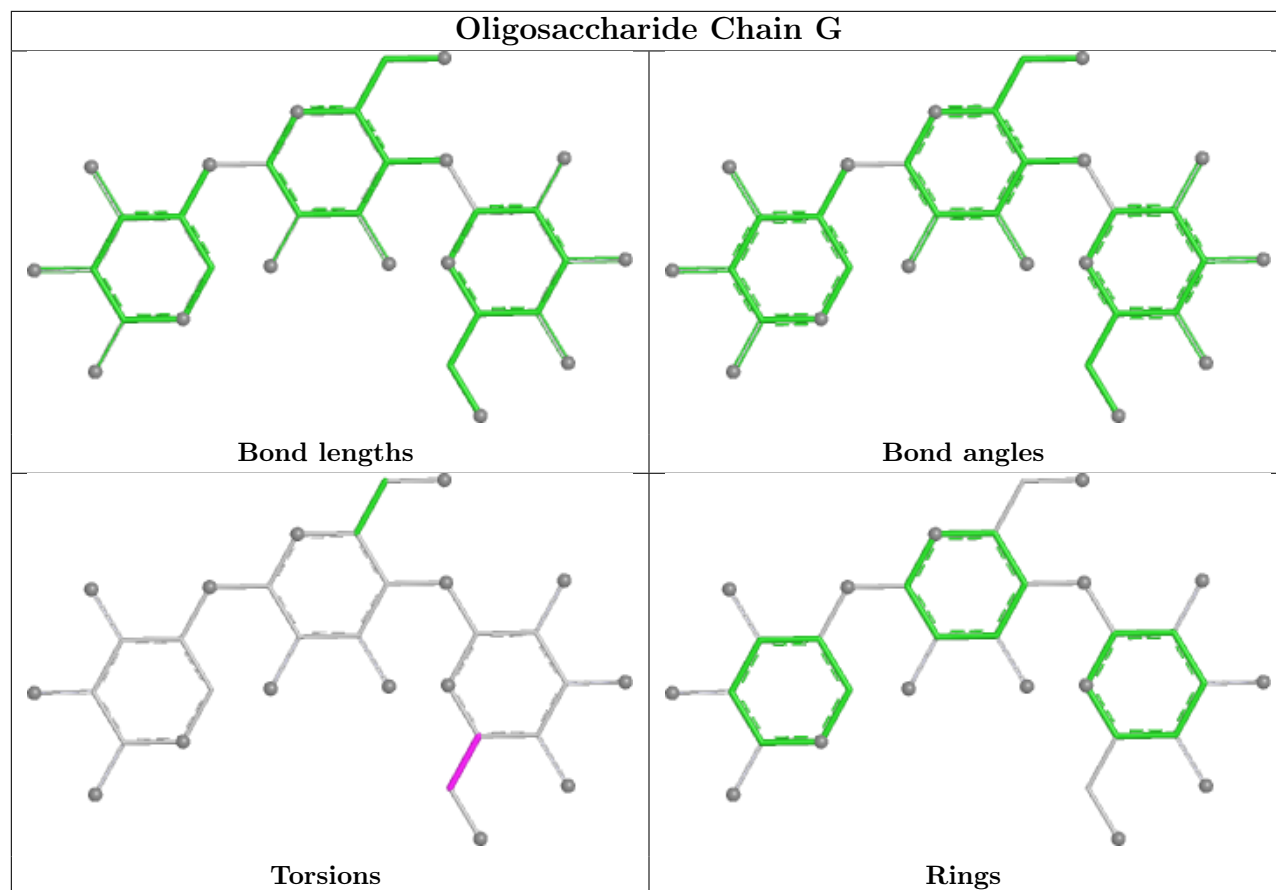
There are no ring outliers.

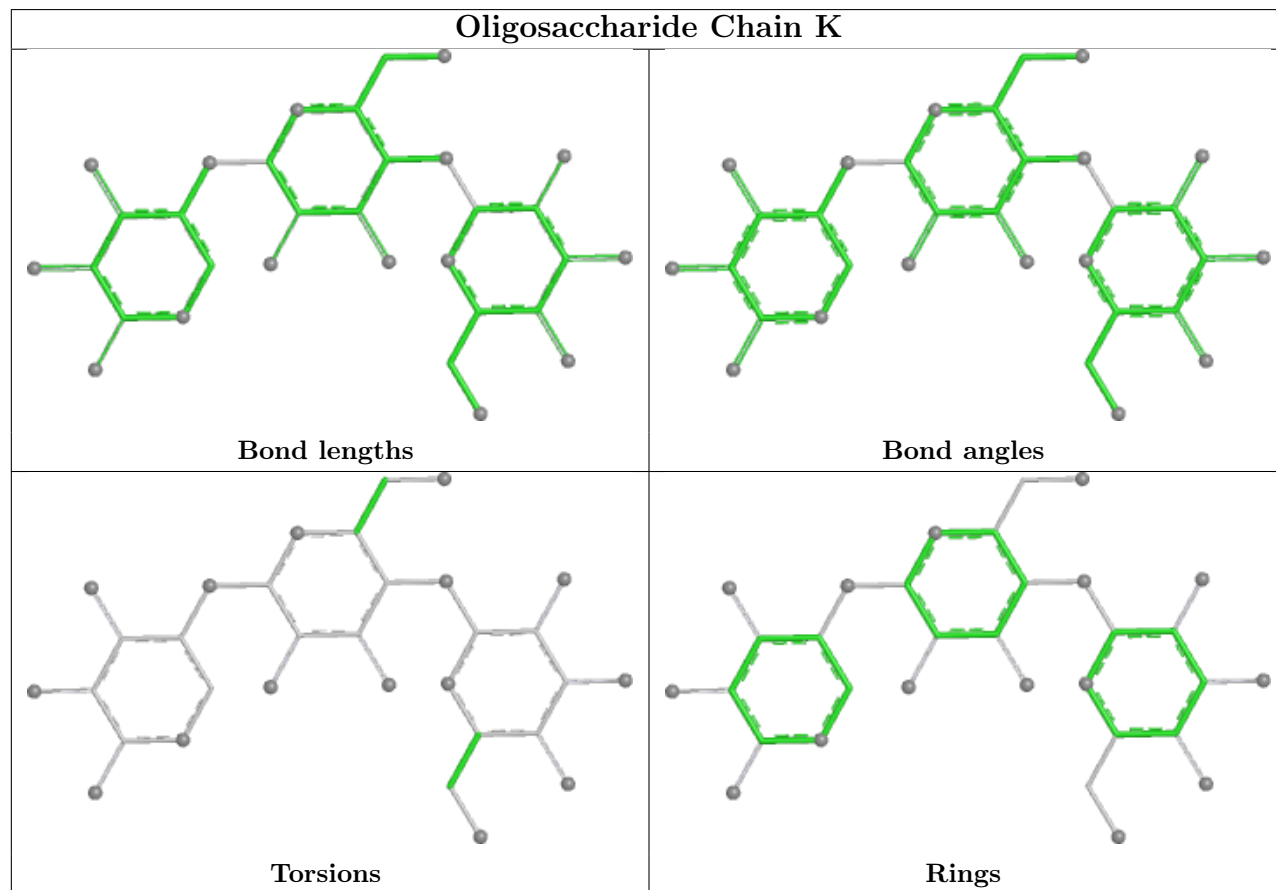
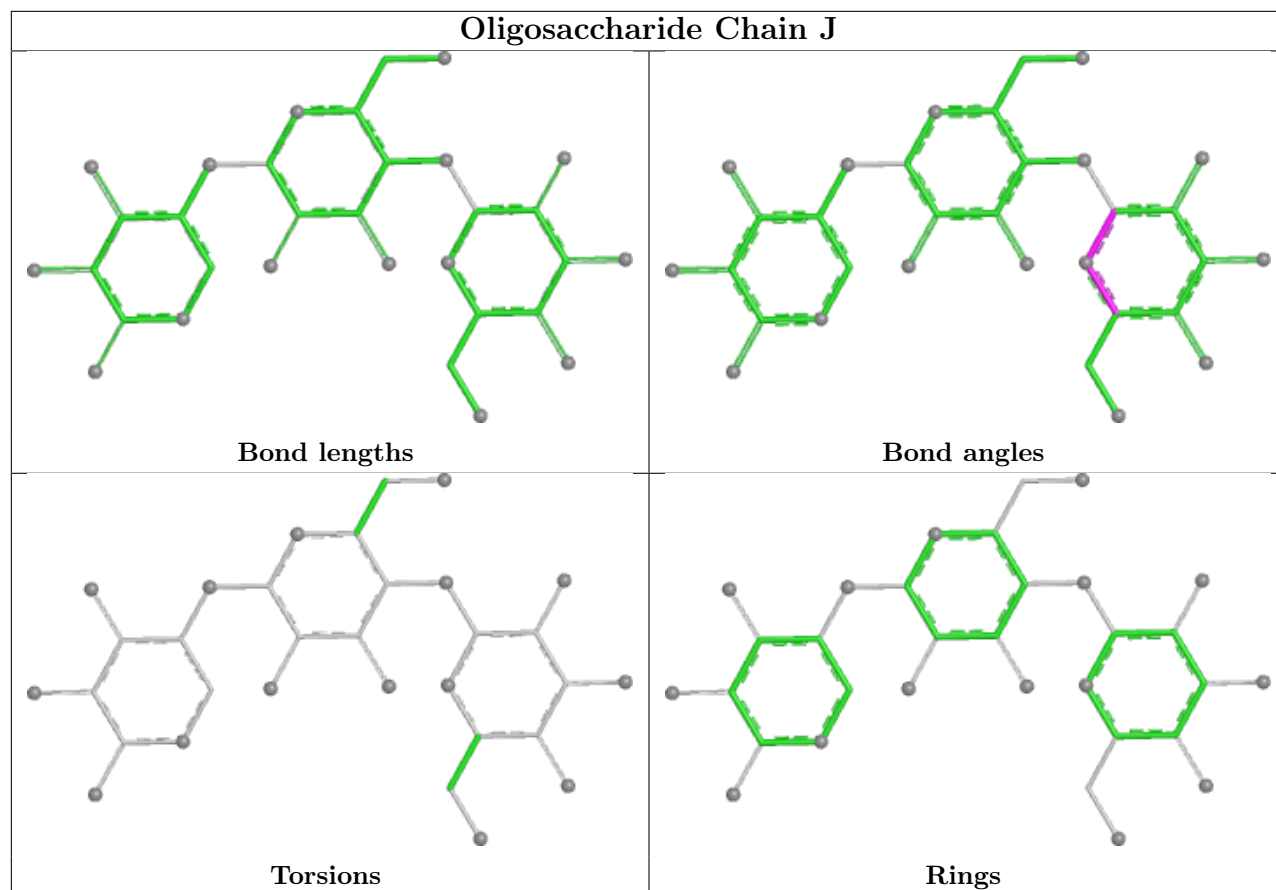
2 monomers are involved in 2 short contacts:

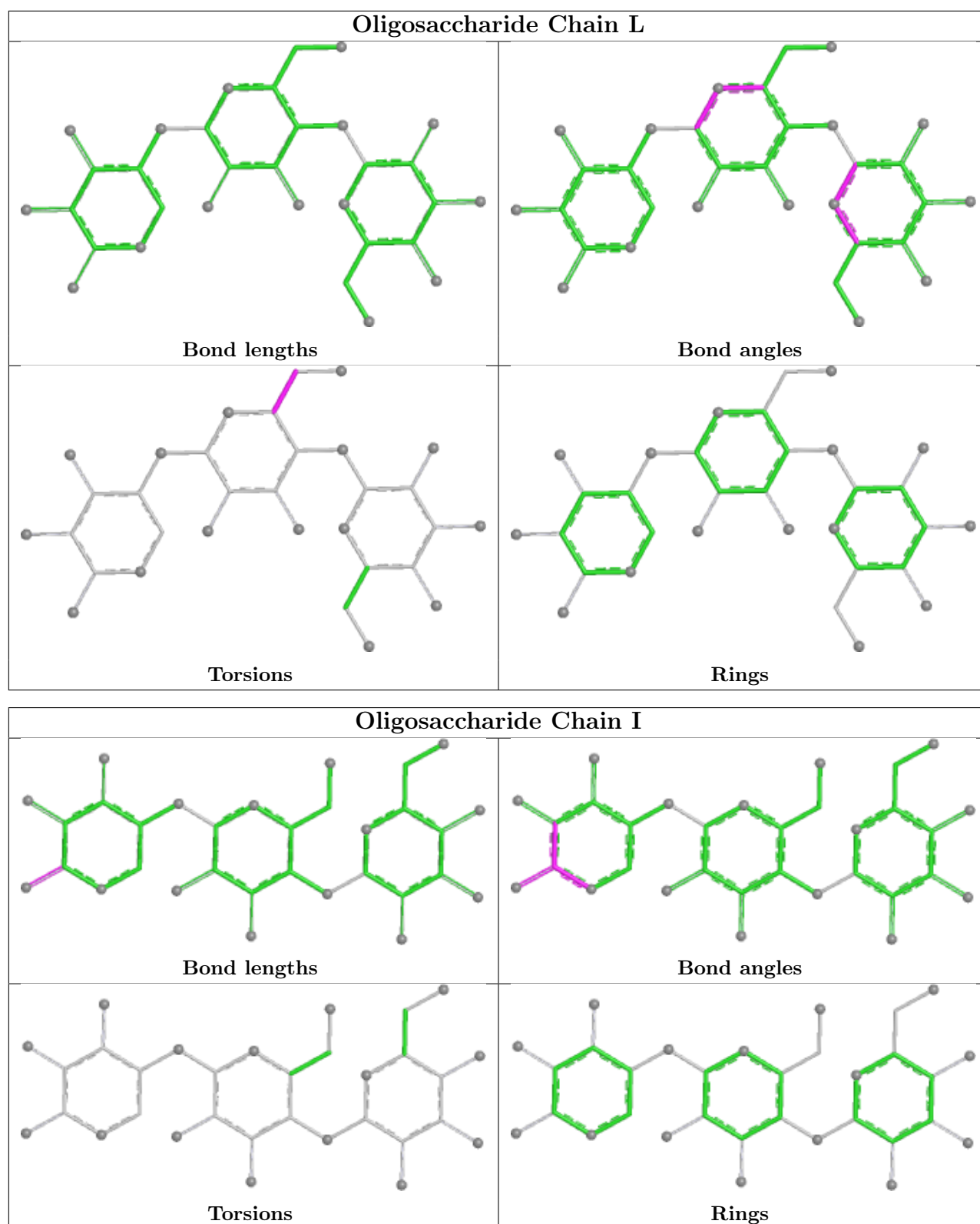
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	I	1	XYS	2	0
3	I	2	GLC	2	0

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









5.6 Ligand geometry [i](#)

Of 4 ligands modelled in this entry, 4 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data [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	516/516 (100%)	-0.80	0 100 100	11, 23, 48, 75	0
1	B	516/516 (100%)	-0.61	3 (0%) 85 87	11, 26, 62, 87	0
1	C	516/516 (100%)	-0.69	1 (0%) 91 92	9, 24, 55, 84	0
1	D	516/516 (100%)	-0.64	2 (0%) 88 90	12, 25, 51, 87	0
All	All	2064/2064 (100%)	-0.68	6 (0%) 90 91	9, 24, 55, 87	0

All (6) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	95	VAL	4.6
1	D	1	ALA	3.2
1	B	95	VAL	2.5
1	B	463	HIS	2.3
1	D	99	CYS	2.3
1	B	432	THR	2.1

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

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

6.3 Carbohydrates [i](#)

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
3	XXS	I	1	10/10	0.69	0.13	0,76,78,80	1

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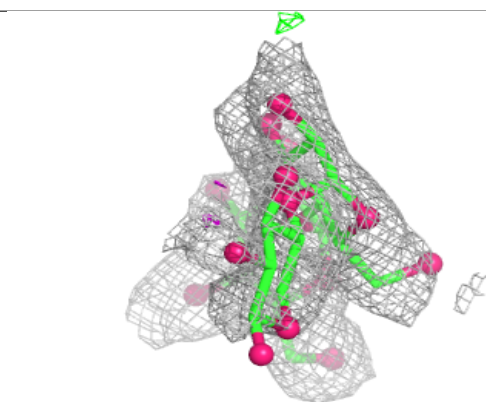
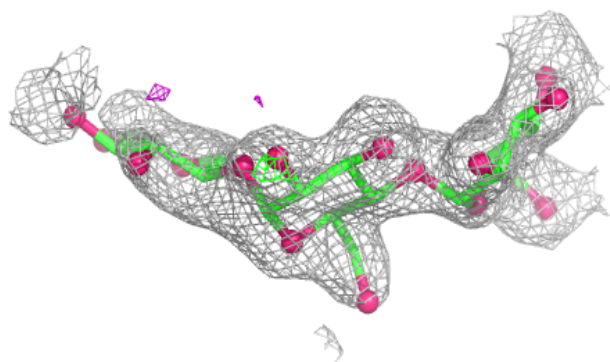
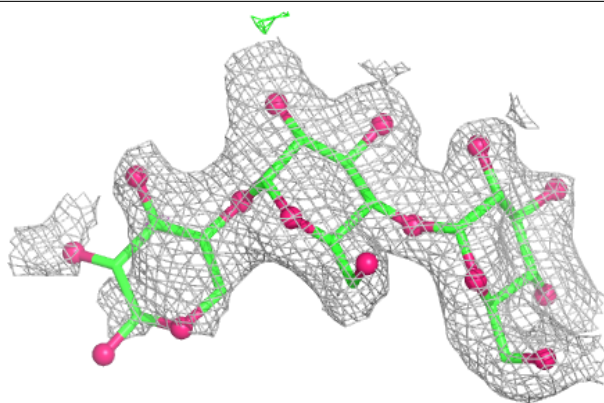
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	XYP	G	1	10/10	0.72	0.12	0,72,75,76	1
2	XYP	L	1	10/10	0.73	0.13	0,80,84,86	1
2	XYP	H	1	10/10	0.73	0.11	0,63,66,69	1
2	XYP	K	1	10/10	0.74	0.11	0,71,76,80	1
2	XYP	F	1	10/10	0.75	0.10	0,56,59,61	1
2	XYP	E	1	10/10	0.76	0.11	0,71,74,77	1
2	GLC	L	2	11/12	0.82	0.10	53,62,63,65	0
2	GLC	L	3	11/12	0.85	0.08	42,45,48,49	0
3	GLC	I	2	11/12	0.85	0.09	40,50,56,60	0
3	GLC	I	3	11/12	0.85	0.09	41,49,51,52	0
2	GLC	G	3	11/12	0.86	0.09	36,48,51,52	0
2	GLC	K	2	11/12	0.86	0.08	46,50,53,55	0
2	XYP	J	1	10/10	0.87	0.10	0,54,56,58	1
2	GLC	K	3	11/12	0.87	0.09	34,42,45,45	0
2	GLC	G	2	11/12	0.89	0.08	43,51,54,55	0
2	GLC	E	2	11/12	0.91	0.07	37,43,48,52	0
2	GLC	F	2	11/12	0.92	0.07	32,34,37,38	0
2	GLC	H	2	11/12	0.92	0.06	35,39,44,46	0
2	GLC	E	3	11/12	0.94	0.07	29,40,44,45	0
2	GLC	J	2	11/12	0.95	0.04	21,27,30,33	0
2	GLC	H	3	11/12	0.95	0.06	24,32,33,34	0
2	GLC	F	3	11/12	0.96	0.05	18,25,29,35	0
2	GLC	J	3	11/12	0.98	0.04	19,22,27,30	0

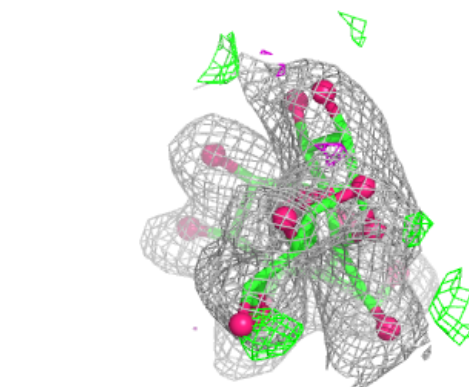
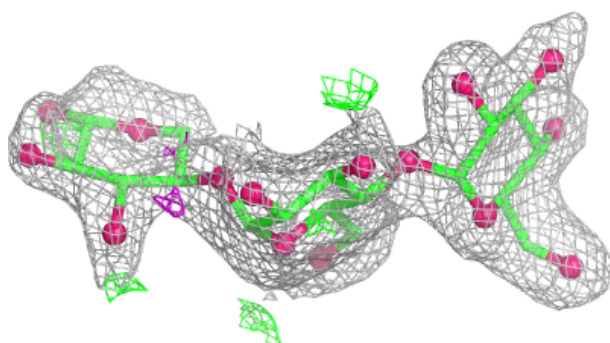
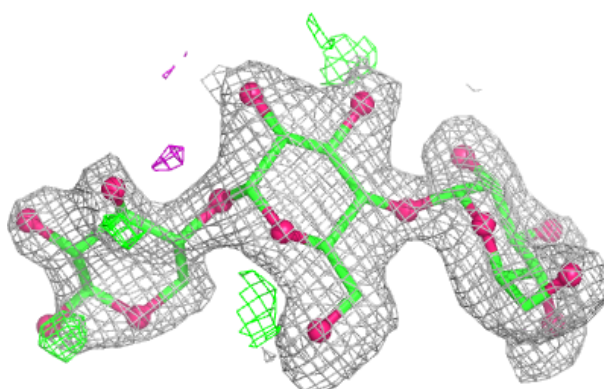
The following is a graphical depiction of the model fit to experimental electron density for oligosaccharide. Each fit is shown from different orientation to approximate a three-dimensional view.

Electron density around Chain E:

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

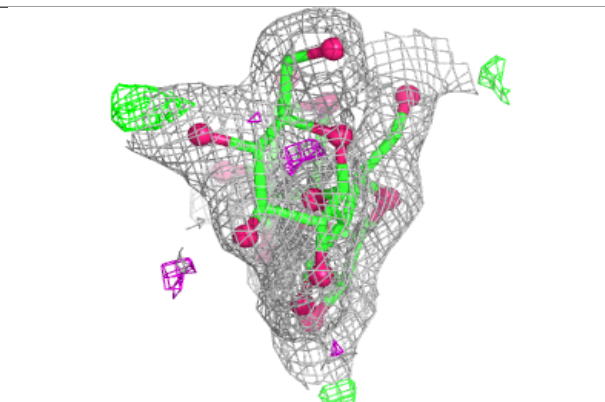
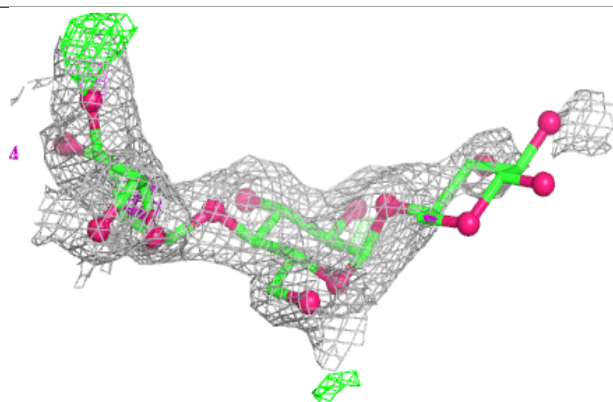
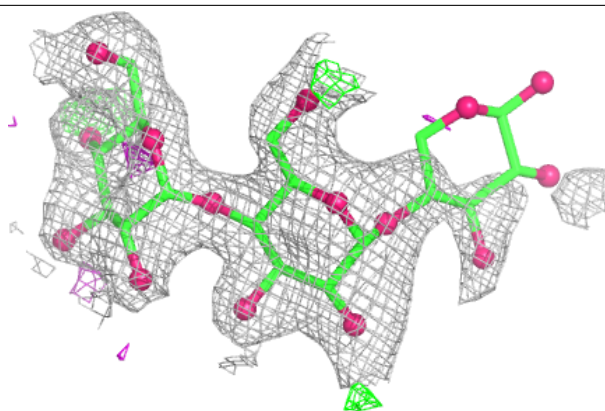
**Electron density around Chain F:**

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

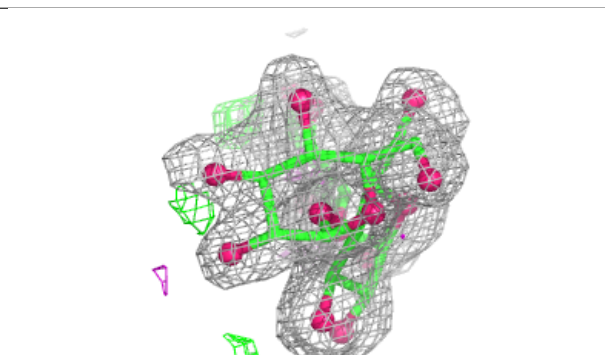
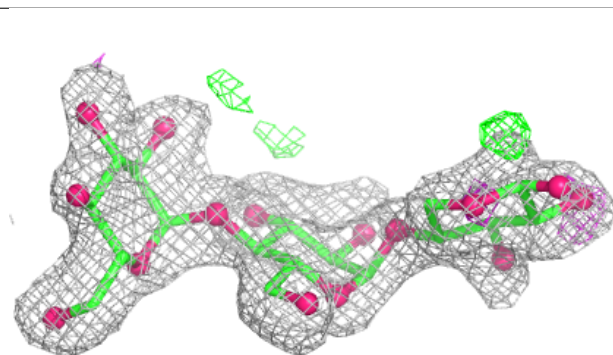
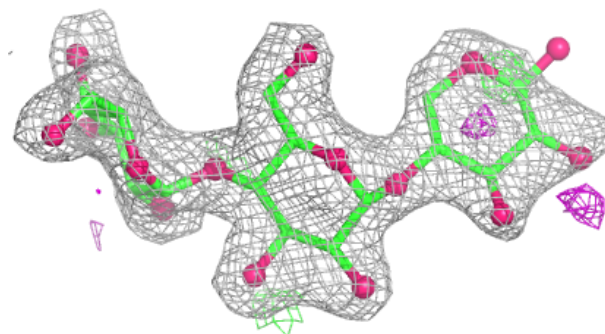


Electron density around Chain G:

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

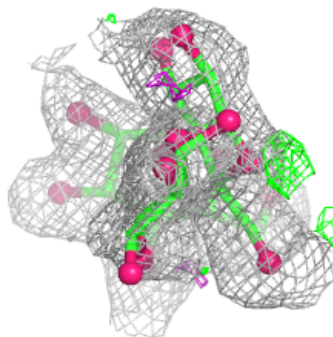
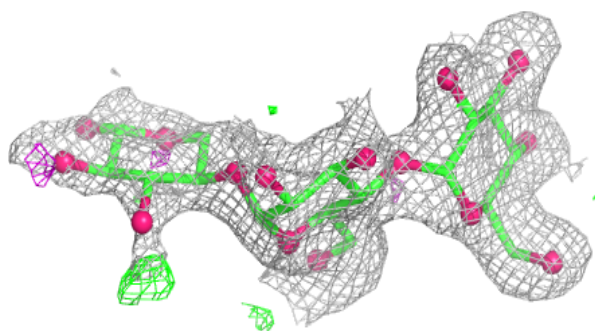
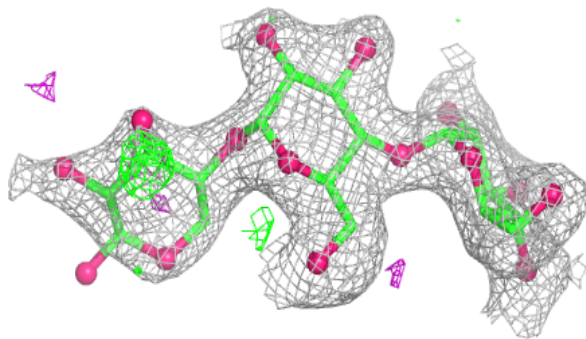
**Electron density around Chain H:**

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

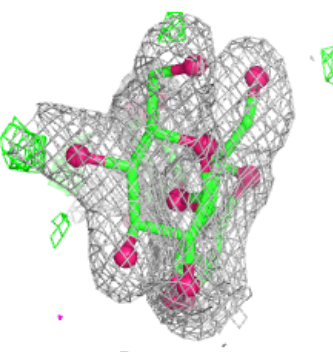
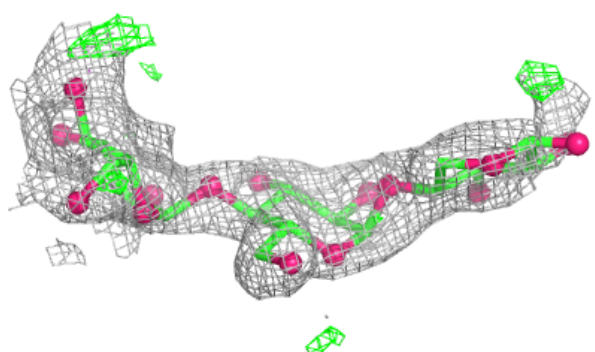
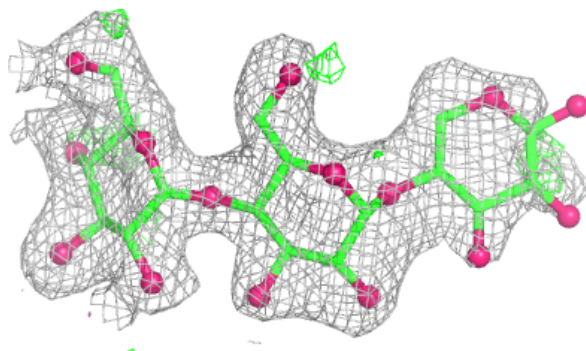


Electron density around Chain J:

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

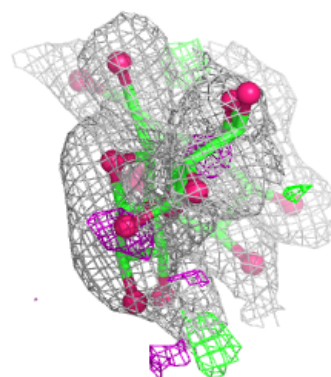
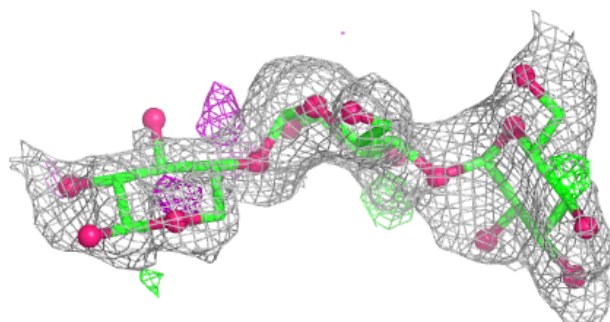
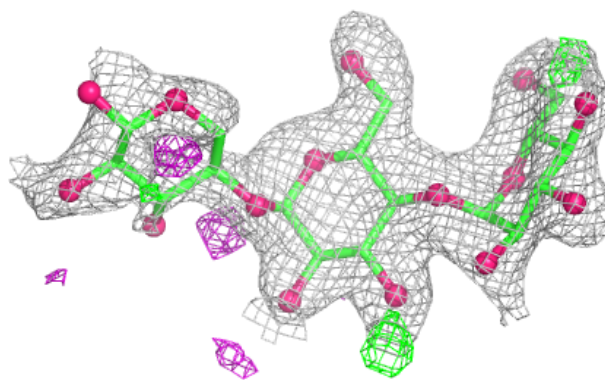
**Electron density around Chain K:**

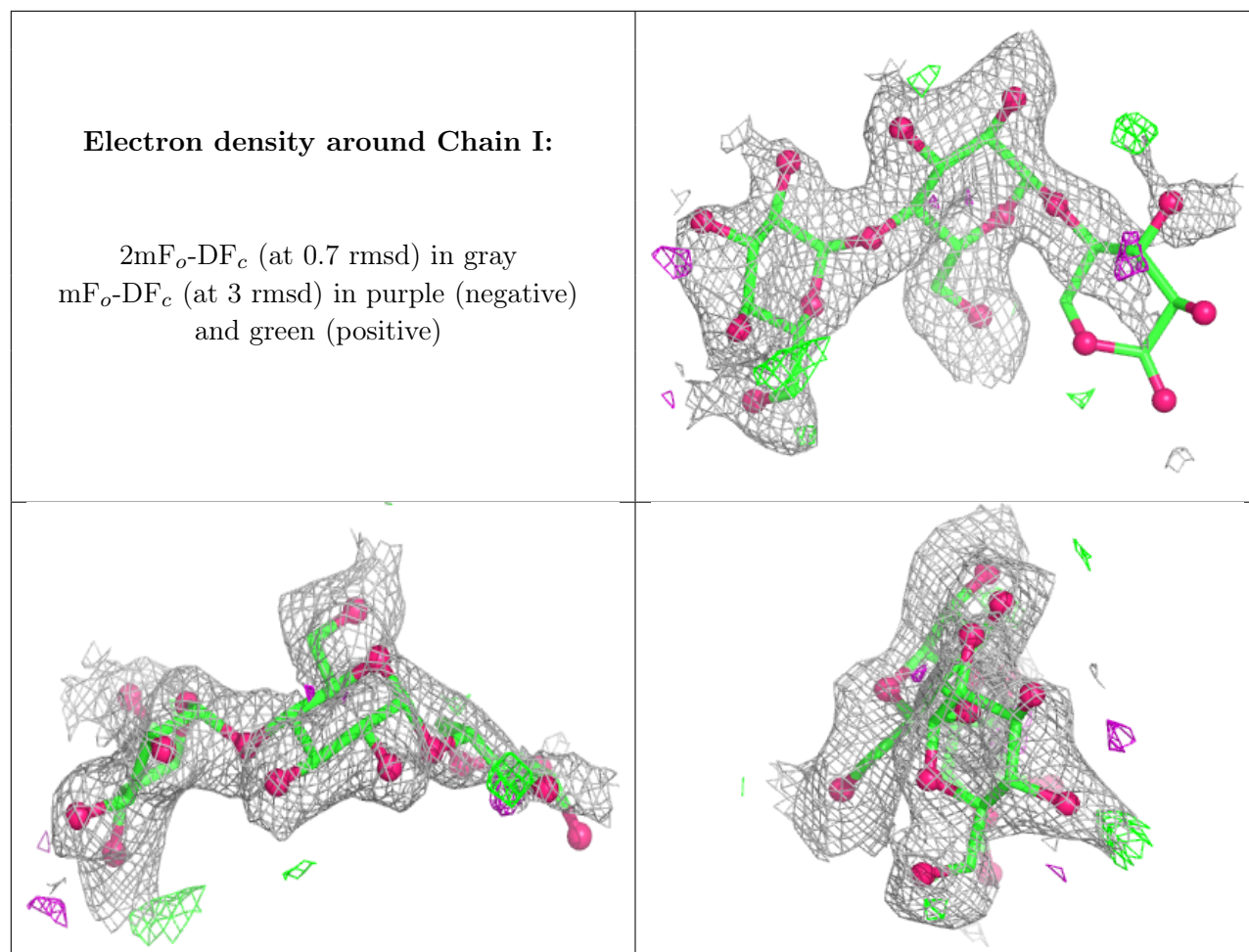
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around Chain L:

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





6.4 Ligands [i](#)

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
4	CA	A	701	1/1	0.99	0.01	18,18,18,18	0
4	CA	C	701	1/1	0.99	0.01	13,13,13,13	0
4	CA	D	701	1/1	0.99	0.01	19,19,19,19	0
4	CA	B	701	1/1	1.00	0.01	15,15,15,15	0

6.5 Other polymers [i](#)

There are no such residues in this entry.