



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 7, 2026 – 01:27 AM UTC

PDB ID : 9M74 / pdb_00009m74
Title : Crystal structure of MBP-fused BIL1/BZR1 (21-104) in complex with double-stranded DNA containing CACAGCTGTG
Authors : Shohei, N.; Masaru, T.; Takuya, M.
Deposited on : 2025-03-09
Resolution : 2.23 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:

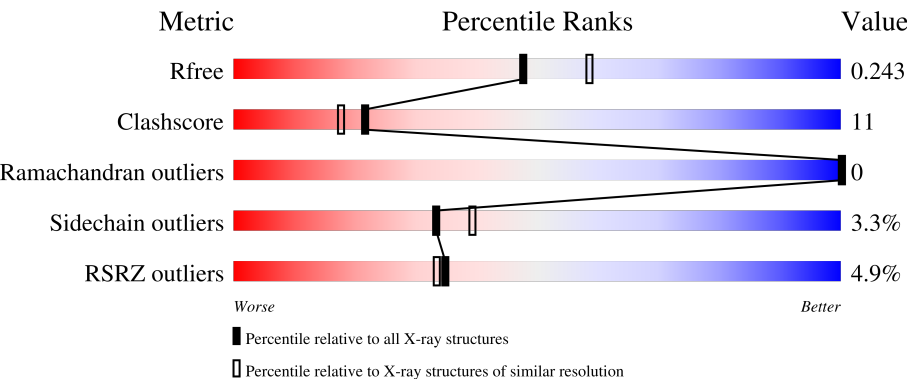
MolProbity : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 2.0
EDS : 3.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (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 2.23 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.



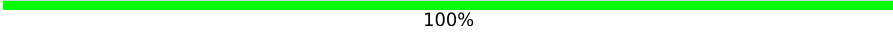
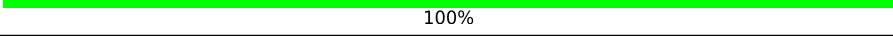
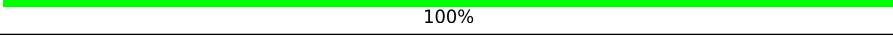
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R _{free}	180053	3416 (2.26-2.22)
Clashscore	190562	3556 (2.26-2.22)
Ramachandran outliers	187476	3500 (2.26-2.22)
Sidechain outliers	187428	3501 (2.26-2.22)
RSRZ outliers	180081	3415 (2.26-2.22)

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	453	3% 83% 12%
1	B	453	2% 83% 13%
1	C	453	2% 75% 19%
1	D	453	12% 68% 25%
2	E	15	80% 20%

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Mol	Chain	Length	Quality of chain
2	F	15	
2	G	15	
2	H	15	
3	J	2	
3	K	2	
3	L	2	
3	M	2	

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 15119 atoms, of which 18 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 Maltodextrin-binding protein, Protein BRASSINAZOLE-RESISTANT 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	C	437	Total	C	N	O	S	0	0	0
			3402	2171	582	640	9			
1	D	436	Total	C	N	O	S	0	0	0
			3394	2166	581	639	8			
1	A	437	Total	C	N	O	S	0	0	0
			3402	2171	582	640	9			
1	B	436	Total	C	N	O	S	0	0	0
			3394	2166	581	639	8			

There are 36 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	-348	MET	-	initiating methionine	UNP A0A4P1LXE0
C	-266	ALA	ASP	engineered mutation	UNP A0A4P1LXE0
C	-265	ALA	LYS	engineered mutation	UNP A0A4P1LXE0
C	-176	ALA	GLU	engineered mutation	UNP A0A4P1LXE0
C	-175	ALA	ASN	engineered mutation	UNP A0A4P1LXE0
C	-109	ALA	LYS	engineered mutation	UNP A0A4P1LXE0
C	11	ALA	GLU	engineered mutation	UNP A0A4P1LXE0
C	14	ALA	LYS	engineered mutation	UNP A0A4P1LXE0
C	15	ALA	ASP	engineered mutation	UNP A0A4P1LXE0
D	-348	MET	-	initiating methionine	UNP A0A4P1LXE0
D	-266	ALA	ASP	engineered mutation	UNP A0A4P1LXE0
D	-265	ALA	LYS	engineered mutation	UNP A0A4P1LXE0
D	-176	ALA	GLU	engineered mutation	UNP A0A4P1LXE0
D	-175	ALA	ASN	engineered mutation	UNP A0A4P1LXE0
D	-109	ALA	LYS	engineered mutation	UNP A0A4P1LXE0
D	11	ALA	GLU	engineered mutation	UNP A0A4P1LXE0
D	14	ALA	LYS	engineered mutation	UNP A0A4P1LXE0
D	15	ALA	ASP	engineered mutation	UNP A0A4P1LXE0
A	-348	MET	-	initiating methionine	UNP A0A4P1LXE0
A	-266	ALA	ASP	engineered mutation	UNP A0A4P1LXE0

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Chain	Residue	Modelled	Actual	Comment	Reference
A	-265	ALA	LYS	engineered mutation	UNP A0A4P1LXE0
A	-176	ALA	GLU	engineered mutation	UNP A0A4P1LXE0
A	-175	ALA	ASN	engineered mutation	UNP A0A4P1LXE0
A	-109	ALA	LYS	engineered mutation	UNP A0A4P1LXE0
A	11	ALA	GLU	engineered mutation	UNP A0A4P1LXE0
A	14	ALA	LYS	engineered mutation	UNP A0A4P1LXE0
A	15	ALA	ASP	engineered mutation	UNP A0A4P1LXE0
B	-348	MET	-	initiating methionine	UNP A0A4P1LXE0
B	-266	ALA	ASP	engineered mutation	UNP A0A4P1LXE0
B	-265	ALA	LYS	engineered mutation	UNP A0A4P1LXE0
B	-176	ALA	GLU	engineered mutation	UNP A0A4P1LXE0
B	-175	ALA	ASN	engineered mutation	UNP A0A4P1LXE0
B	-109	ALA	LYS	engineered mutation	UNP A0A4P1LXE0
B	11	ALA	GLU	engineered mutation	UNP A0A4P1LXE0
B	14	ALA	LYS	engineered mutation	UNP A0A4P1LXE0
B	15	ALA	ASP	engineered mutation	UNP A0A4P1LXE0

- Molecule 2 is a DNA chain called DNA (5'-D(*TP*TP*CP*AP*CP*AP*GP*CP*TP*GP*TP*GP*AP*AP*A)-3').

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	G	15	Total	C	N	O	P	0	0	0
			305	147	57	87	14			
2	H	15	Total	C	N	O	P	0	0	0
			305	147	57	87	14			
2	E	15	Total	C	N	O	P	0	0	0
			305	147	57	87	14			
2	F	15	Total	C	N	O	P	0	0	0
			305	147	57	87	14			

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



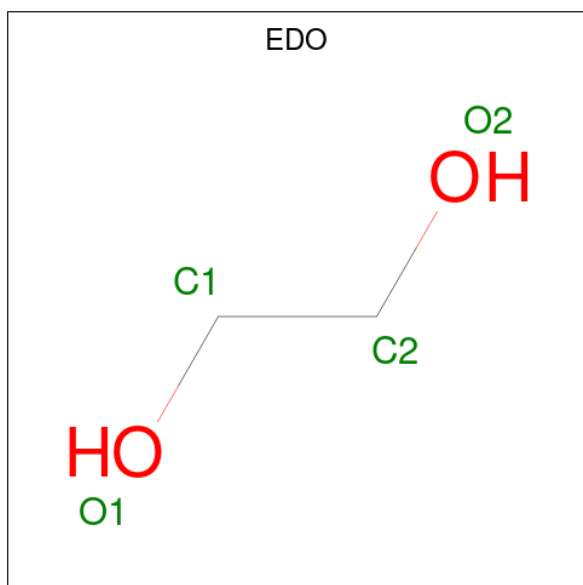
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace
3	J	2	Total	C	O	0	0	0
			23	12	11			
3	K	2	Total	C	O	0	0	0
			23	12	11			

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

- Molecule 4 is 1,2-ETHANEDIOL (CCD ID: EDO) (formula: $C_2H_6O_2$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	C	1	Total	C	H	O	0	0
			10	2	6	2		
4	A	1	Total	C	H	O	0	0
			10	2	6	2		
4	B	1	Total	C	H	O	0	0
			10	2	6	2		

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	C	27	Total	O	0	0
			27	27		
5	D	12	Total	O	0	0
			12	12		
5	G	2	Total	O	0	0
			2	2		
5	H	4	Total	O	0	0
			4	4		

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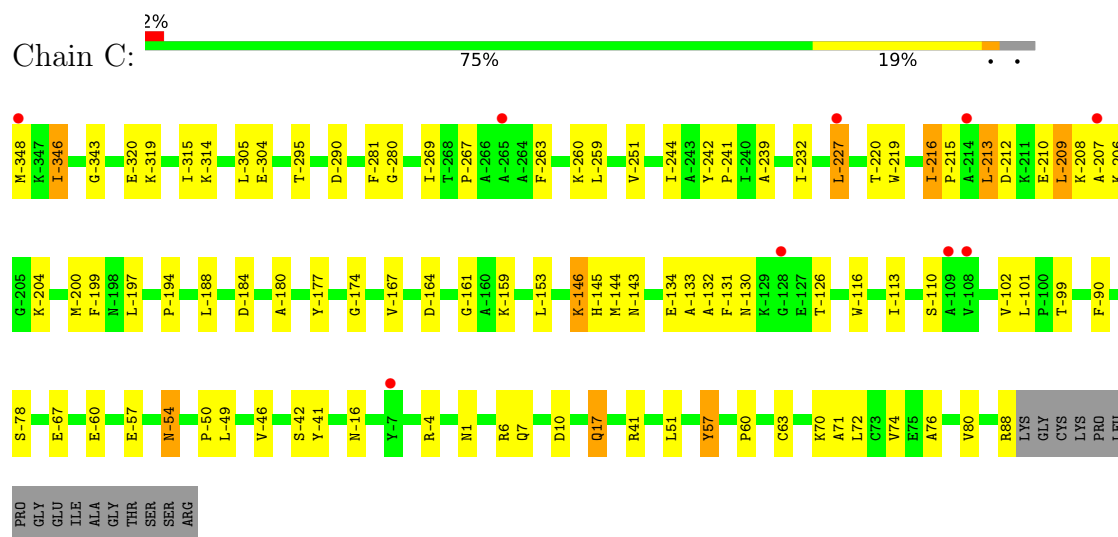
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	49	Total 49	O 49	0	0
5	B	77	Total 77	O 77	0	0
5	E	4	Total 4	O 4	0	0
5	F	10	Total 10	O 10	0	0

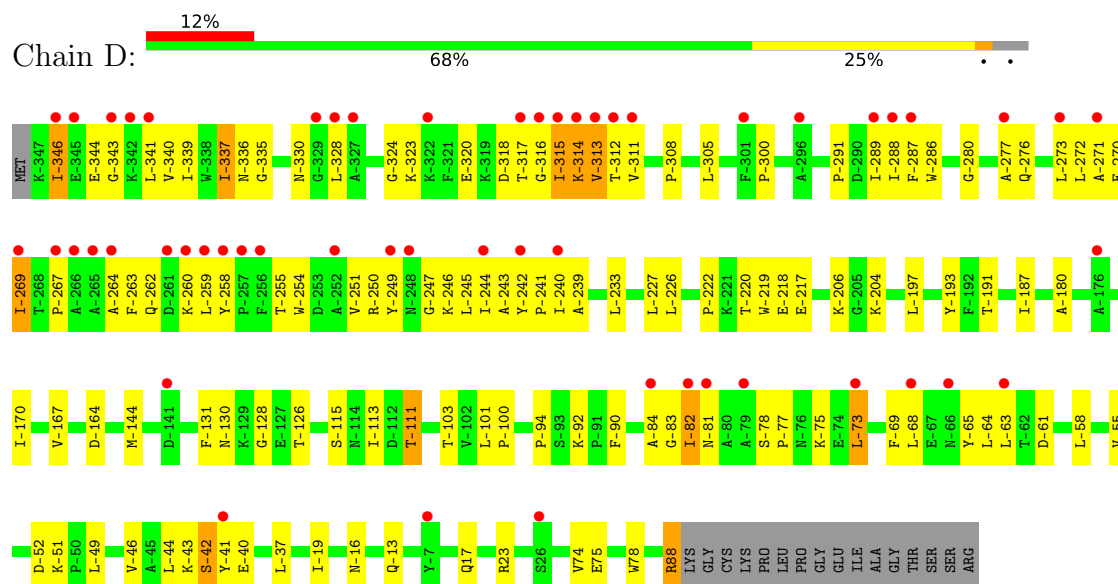
3 Residue-property plots [i](#)

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

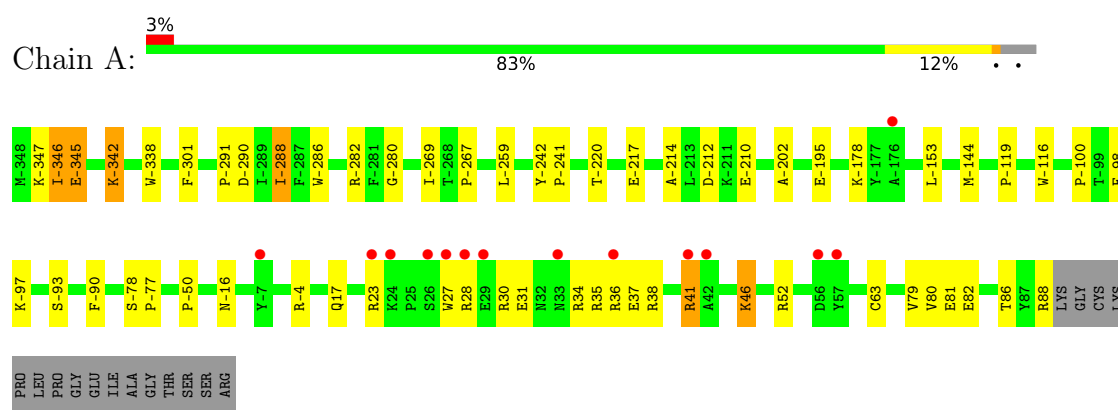
- Molecule 1: Maltodextrin-binding protein,Protein BRASSINAZOLE-RESISTANT 1



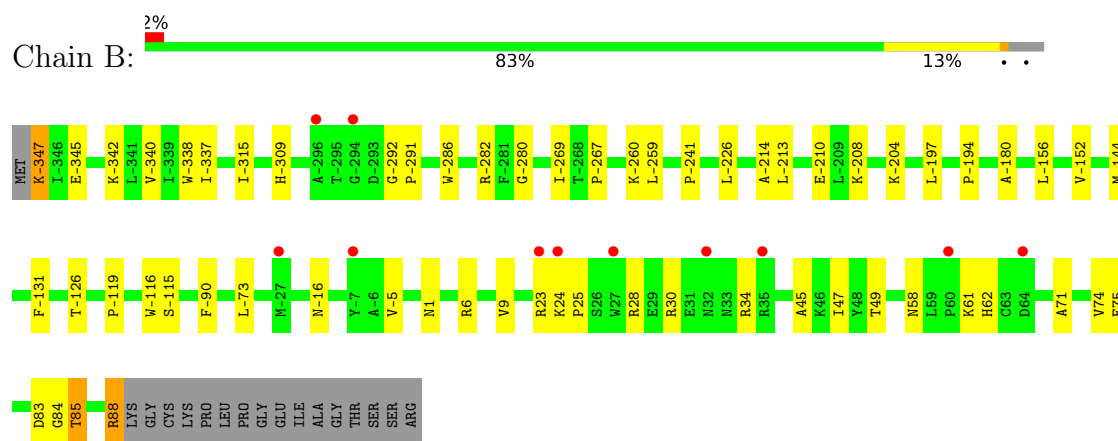
- Molecule 1: Maltodextrin-binding protein,Protein BRASSINAZOLE-RESISTANT 1



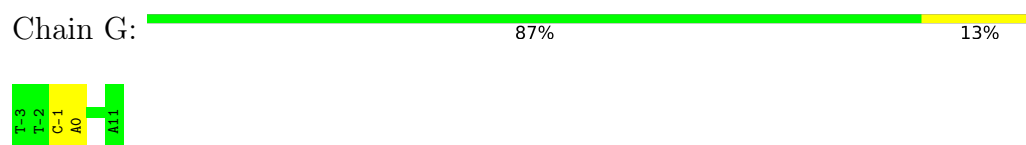
- Molecule 1: Maltodextrin-binding protein,Protein BRASSINAZOLE-RESISTANT 1



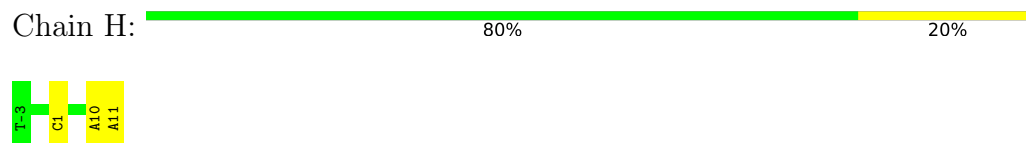
- Molecule 1: Maltodextrin-binding protein, Protein BRASSINAZOLE-RESISTANT 1



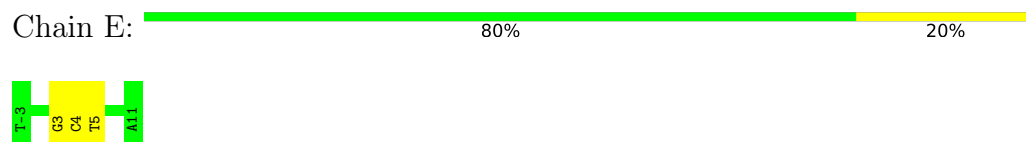
- Molecule 2: DNA (5'-D(*TP*TP*CP*AP*CP*AP*GP*CP*TP*GP*TP*GP*AP*AP*A)-3')



- Molecule 2: DNA (5'-D(*TP*TP*CP*AP*CP*AP*GP*CP*TP*GP*TP*GP*AP*AP*A)-3')



- Molecule 2: DNA (5'-D(*TP*TP*CP*AP*CP*AP*GP*CP*TP*GP*TP*GP*AP*AP*A)-3')

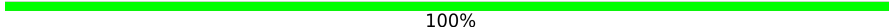


- Molecule 2: DNA (5'-D(*TP*TP*CP*AP*CP*AP*GP*CP*TP*GP*TP*GP*AP*AP*A)-3')

Chain F: 



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

Chain J: 

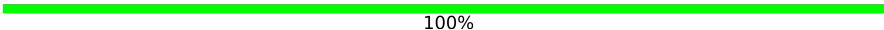


- Molecule 3: alpha-D-glucopyranose-(1-4)-alpha-D-glucopyranose

Chain K: 



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

Chain L: 



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

Chain M: 



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	102.84Å 92.51Å 111.74Å 90.00° 100.17° 90.00°	Depositor
Resolution (Å)	43.58 – 2.23 43.58 – 2.23	Depositor EDS
% Data completeness (in resolution range)	99.9 (43.58-2.23) 99.9 (43.58-2.23)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.96 (at 2.22Å)	Xtriage
Refinement program	PHENIX 1.21.2_5419	Depositor
R, R_{free}	0.215 , 0.244 0.214 , 0.243	Depositor DCC
R_{free} test set	5006 reflections (4.98%)	wwPDB-VP
Wilson B-factor (Å ²)	39.6	Xtriage
Anisotropy	0.374	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 36.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	15119	wwPDB-VP
Average B, all atoms (Å ²)	54.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 44.00 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 1.6261e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

¹Intensities estimated from amplitudes.

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

5 Model quality [i](#)

5.1 Standard geometry [i](#)

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

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

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z > 5$	RMSZ	$\# Z > 5$
1	A	0.29	0/3481	0.36	0/4723
1	B	0.40	2/3473 (0.1%)	0.40	0/4713
1	C	0.33	1/3481 (0.0%)	0.39	2/4723 (0.0%)
1	D	0.37	0/3473	0.46	1/4713 (0.0%)
2	E	0.23	0/342	0.51	0/526
2	F	0.22	0/342	0.46	0/526
2	G	0.24	0/342	0.48	0/526
2	H	0.25	0/342	0.49	0/526
All	All	0.34	3/15276 (0.0%)	0.41	3/20976 (0.0%)

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	-180	ALA	C-O	-7.75	1.17	1.23
1	C	-54	ASN	C-O	-5.17	1.18	1.24
1	B	-292	GLY	C-O	-5.13	1.17	1.24

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	-146	LYS	N-CA-C	5.49	119.00	111.54
1	D	-335	GLY	N-CA-C	5.44	120.44	114.40
1	C	57	TYR	N-CA-C	5.42	119.52	113.02

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	3402	0	3374	57	0
1	B	3394	0	3365	41	0
1	C	3402	0	3374	74	0
1	D	3394	0	3365	144	0
2	E	305	0	171	2	0
2	F	305	0	171	4	0
2	G	305	0	171	1	0
2	H	305	0	171	2	0
3	J	23	0	21	0	0
3	K	23	0	21	0	0
3	L	23	0	21	0	0
3	M	23	0	21	1	0
4	A	4	6	6	0	0
4	B	4	6	6	0	0
4	C	4	6	6	0	0
5	A	49	0	0	0	0
5	B	77	0	0	0	0
5	C	27	0	0	0	0
5	D	12	0	0	1	0
5	E	4	0	0	0	0
5	F	10	0	0	0	0
5	G	2	0	0	0	0
5	H	4	0	0	2	0
All	All	15101	18	14264	316	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 11.

All (316) 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:D:-320:GLU:HG2	1:D:-315:ILE:O	1.36	1.21
1:D:-315:ILE:HD12	1:D:-73:LEU:HD11	1.19	1.17
1:D:-263:PHE:HA	1:D:-260:LYS:CD	1.80	1.11
1:D:-263:PHE:HA	1:D:-260:LYS:HD3	1.28	1.06

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:-346:ILE:HG22	1:A:-342:LYS:NZ	1.70	1.05
1:D:-41:TYR:CE2	1:D:-37:LEU:HD11	1.92	1.04
1:D:-315:ILE:CD1	1:D:-73:LEU:HD21	1.90	1.01
1:D:-315:ILE:HD12	1:D:-73:LEU:CD1	1.92	0.99
1:B:88:ARG:HH11	1:B:88:ARG:HB3	1.25	0.99
1:D:-315:ILE:CD1	1:D:-73:LEU:HD11	1.97	0.94
1:D:-337:ILE:HG13	1:D:-287:PHE:HB2	1.50	0.94
1:D:-191:THR:HG22	1:D:-187:ILE:HD11	1.49	0.92
1:C:-215:PRO:HA	1:C:-145:HIS:CD2	2.05	0.92
1:D:-191:THR:HG22	1:D:-187:ILE:CD1	2.00	0.91
1:C:-269:ILE:HG22	1:C:-267:PRO:HD3	1.50	0.91
1:D:-249:TYR:HB3	1:D:-244:ILE:HD11	1.51	0.90
1:D:-264:ALA:O	1:D:-260:LYS:HD2	1.73	0.89
1:D:-250:ARG:HG2	1:D:-245:LEU:HG	1.55	0.89
1:A:-346:ILE:HG22	1:A:-342:LYS:HZ3	1.36	0.88
1:B:30:ARG:HE	1:B:34:ARG:HH22	1.20	0.87
1:A:-346:ILE:HD12	1:A:-77:PRO:HD3	1.57	0.86
1:D:-250:ARG:CG	1:D:-245:LEU:HG	2.08	0.83
1:D:-315:ILE:HD12	1:D:-73:LEU:HD21	1.59	0.82
1:A:-346:ILE:HG22	1:A:-342:LYS:HZ1	1.43	0.81
1:D:-315:ILE:HD11	1:D:-73:LEU:HD21	1.60	0.80
1:D:-270:GLU:OE1	1:D:-246:LYS:HE2	1.81	0.79
1:D:-346:ILE:HD13	1:D:-77:PRO:HD3	1.64	0.79
1:A:-220:THR:HG22	1:A:-217:GLU:OE1	1.83	0.78
1:D:88:ARG:HH11	1:D:88:ARG:HG3	1.48	0.78
1:D:-320:GLU:CD	1:D:-314:LYS:HA	2.08	0.78
1:D:-263:PHE:HA	1:D:-260:LYS:HD2	1.66	0.77
1:C:-54:ASN:ND2	1:C:-49:LEU:H	1.82	0.77
1:D:-280:GLY:HA3	1:D:-16:ASN:O	1.85	0.76
1:D:-343:GLY:O	1:D:-315:ILE:HG23	1.86	0.75
1:C:-346:ILE:HD13	1:C:-346:ILE:H	1.51	0.73
1:D:-264:ALA:C	1:D:-260:LYS:HD2	2.13	0.73
1:A:-259:LEU:HD23	1:A:-241:PRO:HG2	1.69	0.73
1:C:51:LEU:HD11	1:C:72:LEU:HD22	1.69	0.73
1:D:-19:ILE:HD12	1:D:-19:ILE:O	1.89	0.72
1:B:-269:ILE:HG22	1:B:-267:PRO:HD3	1.70	0.72
1:D:-337:ILE:HG13	1:D:-287:PHE:CB	2.19	0.72
1:D:-315:ILE:HD12	1:D:-73:LEU:CD2	2.19	0.72
1:D:-320:GLU:OE1	1:D:-314:LYS:HA	1.88	0.72
1:C:-209:LEU:HD12	1:C:-204:LYS:HB2	1.73	0.71
1:D:-315:ILE:CD1	1:D:-73:LEU:CD2	2.68	0.71

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:-344:GLU:HG2	1:D:-77:PRO:HB2	1.72	0.70
1:A:31:GLU:O	1:A:35:ARG:HG2	1.91	0.70
1:A:37:GLU:O	1:A:41:ARG:HG3	1.91	0.69
1:C:-212:ASP:HB3	1:C:-145:HIS:CD2	2.28	0.69
1:C:-212:ASP:HB3	1:C:-145:HIS:HD2	1.56	0.69
1:C:41:ARG:HD3	5:H:101:HOH:O	1.92	0.69
1:D:-64:LEU:O	1:D:-58:LEU:CD2	2.41	0.69
1:C:-280:GLY:HA3	1:C:-16:ASN:O	1.93	0.68
1:D:-41:TYR:CD2	1:D:-37:LEU:HD11	2.29	0.68
1:A:-345:GLU:H	1:A:-342:LYS:CE	2.07	0.67
1:A:-342:LYS:HD2	1:A:-342:LYS:C	2.19	0.67
1:A:81:GLU:CD	1:A:82:GLU:H	2.02	0.67
1:C:-281:PHE:HB3	1:C:-244:ILE:HD13	1.76	0.67
1:A:-269:ILE:HG22	1:A:-267:PRO:HD3	1.77	0.67
1:B:88:ARG:HH11	1:B:88:ARG:CB	2.03	0.66
1:D:-78:SER:O	1:D:-75:LYS:NZ	2.27	0.66
1:A:-345:GLU:O	1:A:-342:LYS:HE3	1.95	0.66
1:A:-346:ILE:HD13	1:A:-78:SER:HA	1.78	0.65
1:D:-315:ILE:HD12	1:D:-73:LEU:CG	2.26	0.65
1:B:24:LYS:HD3	1:B:25:PRO:N	2.10	0.65
1:B:-156:LEU:O	1:B:-152:VAL:HG23	1.96	0.64
1:D:-269:ILE:HG23	1:D:-267:PRO:HD3	1.79	0.64
1:D:-324:GLY:O	1:D:-313:VAL:HG21	1.98	0.64
1:B:45:ALA:O	1:B:49:THR:HG23	1.97	0.64
1:D:-344:GLU:CG	1:D:-77:PRO:HB2	2.27	0.64
1:A:-153:LEU:CD1	1:A:-144:MET:HE1	2.27	0.64
1:D:-324:GLY:O	1:D:-313:VAL:CG2	2.46	0.64
1:D:-240:ILE:HD13	1:D:-64:LEU:HD21	1.80	0.64
1:B:-280:GLY:HA3	1:B:-16:ASN:O	1.98	0.63
1:B:-197:LEU:HD11	1:B:-144:MET:HE3	1.80	0.63
1:D:-259:LEU:HD22	1:D:-44:LEU:HA	1.81	0.63
1:D:-336:ASN:OD1	1:D:-286:TRP:HZ3	1.81	0.62
1:D:-320:GLU:OE1	1:D:-314:LYS:CA	2.48	0.62
1:B:1:ASN:OD1	1:B:6:ARG:NH1	2.33	0.62
1:D:-339:ILE:HG12	1:D:-289:ILE:HD12	1.81	0.61
1:D:-220:THR:HG22	1:D:-217:GLU:OE1	2.00	0.61
1:D:-340:VAL:HG23	1:D:-312:THR:OG1	1.99	0.61
1:D:-341:LEU:CB	1:D:-313:VAL:HG13	2.30	0.61
1:A:31:GLU:OE1	1:A:35:ARG:NH1	2.33	0.61
1:D:-328:LEU:O	1:D:-328:LEU:HD23	2.01	0.61
1:D:-242:TYR:CD2	1:D:-68:LEU:HB3	2.35	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:-115:SER:O	1:D:-111:THR:HG23	2.00	0.61
1:A:41:ARG:NH1	2:E:3:DG:OP2	2.34	0.61
1:D:-242:TYR:CD2	1:D:-68:LEU:HD23	2.36	0.61
1:D:-55:VAL:HG11	1:D:-49:LEU:HD21	1.82	0.60
1:A:46:LYS:NZ	1:B:85:THR:HG22	2.16	0.60
1:D:-52:ASP:C	1:D:-51:LYS:HD2	2.27	0.60
1:B:24:LYS:HD3	1:B:24:LYS:C	2.27	0.60
1:D:-242:TYR:CE2	1:D:-68:LEU:HB3	2.37	0.60
1:D:-81:ASN:HB3	1:D:-78:SER:HB2	1.84	0.60
1:A:-153:LEU:HD11	1:A:-144:MET:HE1	1.83	0.60
1:D:-318:ASP:OD2	1:D:-65:TYR:OH	2.20	0.59
1:D:-64:LEU:O	1:D:-58:LEU:HD21	2.01	0.59
1:A:-301:PHE:CG	1:A:-288:ILE:HD12	2.37	0.59
1:C:-54:ASN:HD22	1:C:-49:LEU:H	1.48	0.59
1:C:-208:LYS:HA	1:C:-204:LYS:O	2.02	0.59
1:D:-206:LYS:NZ	1:D:-206:LYS:HB3	2.17	0.59
1:C:-200:MET:HE2	1:C:-126:THR:HB	1.85	0.58
1:D:-251:VAL:HG22	1:D:-251:VAL:O	2.02	0.58
1:D:-260:LYS:C	1:D:-259:LEU:HD23	2.28	0.58
1:A:-280:GLY:HA3	1:A:-16:ASN:O	2.03	0.58
1:B:-152:VAL:HG21	1:B:9:VAL:CG2	2.34	0.58
1:B:-345:GLU:OE2	1:B:-342:LYS:HE2	2.04	0.57
1:B:24:LYS:NZ	1:B:25:PRO:O	2.31	0.57
1:D:-251:VAL:O	1:D:-251:VAL:HG13	2.03	0.57
1:D:88:ARG:HG3	1:D:88:ARG:NH1	2.15	0.57
1:D:-249:TYR:CB	1:D:-244:ILE:HD11	2.29	0.57
1:D:-61:ASP:OD2	1:D:-42:SER:HB2	2.05	0.57
1:A:-342:LYS:NZ	1:A:-290:ASP:OD2	2.37	0.57
1:C:-199:PHE:O	1:C:-197:LEU:HD13	2.05	0.57
1:B:-226:LEU:HD21	1:B:-213:LEU:HD21	1.87	0.56
1:A:86:THR:OG1	1:B:47:ILE:HA	2.05	0.56
1:D:-246:LYS:HE3	1:D:-246:LYS:HA	1.88	0.56
1:C:51:LEU:CD1	1:C:72:LEU:HD22	2.36	0.56
1:A:-220:THR:HG23	1:A:-217:GLU:H	1.71	0.56
1:B:-208:LYS:HD3	1:B:-204:LYS:O	2.06	0.56
1:C:-212:ASP:OD1	1:C:-208:LYS:HE3	2.05	0.56
1:D:-315:ILE:CD1	1:D:-73:LEU:CG	2.84	0.56
1:A:-345:GLU:H	1:A:-342:LYS:HE3	1.69	0.56
1:D:-258:TYR:CZ	1:D:-43:LYS:HG2	2.40	0.56
1:C:-210:GLU:O	1:C:-207:ALA:HB3	2.05	0.56
1:C:-263:PHE:HA	1:C:-260:LYS:HD2	1.87	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:-346:ILE:CD1	1:A:-77:PRO:HD3	2.31	0.55
1:D:-40:GLU:OE2	1:D:-40:GLU:HA	2.06	0.55
1:D:-276:GLN:HB2	1:D:-249:TYR:HE2	1.71	0.55
1:D:-167:VAL:HG23	1:D:17:GLN:OE1	2.07	0.55
1:D:-269:ILE:HG12	1:D:-242:TYR:CE1	2.42	0.55
1:D:-263:PHE:CA	1:D:-260:LYS:HD2	2.36	0.55
1:D:-330:ASN:HB2	1:D:-52:ASP:OD2	2.06	0.55
1:D:-341:LEU:HB3	1:D:-313:VAL:HG13	1.89	0.55
1:D:-240:ILE:HD13	1:D:-64:LEU:CD2	2.36	0.55
1:A:-269:ILE:HD12	1:A:-242:TYR:CE2	2.42	0.54
1:C:-227:LEU:HD21	1:C:-204:LYS:HG3	1.88	0.54
1:D:-320:GLU:O	1:D:-316:GLY:N	2.37	0.54
1:C:-164:ASP:O	1:C:-159:LYS:HE3	2.08	0.54
1:D:-94:PRO:HG2	1:D:-92:LYS:HE3	1.90	0.54
1:A:88:ARG:NH2	1:B:75:GLU:OE2	2.39	0.54
1:D:-69:PHE:HA	1:D:-65:TYR:HB2	1.88	0.54
1:B:-259:LEU:HD23	1:B:-241:PRO:HG2	1.90	0.54
1:D:-191:THR:CG2	1:D:-187:ILE:HD11	2.30	0.54
1:D:-187:ILE:N	1:D:-187:ILE:HD12	2.23	0.54
1:D:-220:THR:HG22	1:D:-217:GLU:CD	2.33	0.54
1:D:-226:LEU:HD21	1:D:-222:PRO:HD3	1.90	0.53
1:B:71:ALA:O	1:B:74:VAL:HG22	2.08	0.53
1:D:-315:ILE:HD11	1:D:-73:LEU:CD2	2.34	0.53
1:A:-286:TRP:CD1	1:A:-282:ARG:HG3	2.42	0.53
1:B:88:ARG:HB3	1:B:88:ARG:NH1	2.09	0.53
1:C:-200:MET:HE1	1:C:-132:ALA:HB1	1.90	0.52
1:D:-341:LEU:HB2	1:D:-313:VAL:HG13	1.91	0.52
1:D:-164:ASP:HB2	1:D:17:GLN:HB2	1.89	0.52
1:B:-214:ALA:O	1:B:-210:GLU:HG2	2.09	0.52
1:C:-263:PHE:CE2	1:C:-67:GLU:HG2	2.45	0.52
1:D:-239:ALA:HA	1:D:-46:VAL:HA	1.90	0.52
1:D:-263:PHE:CA	1:D:-260:LYS:CD	2.72	0.52
1:C:-346:ILE:HG13	1:C:-78:SER:HA	1.91	0.52
1:D:-64:LEU:O	1:D:-58:LEU:HD23	2.08	0.52
1:B:-347:LYS:HD3	1:B:-340:VAL:HG11	1.92	0.52
1:D:-323:LYS:O	1:D:-320:GLU:HB2	2.10	0.51
1:D:-277:ALA:HB2	1:D:-244:ILE:HG21	1.93	0.51
1:D:-41:TYR:CD2	1:D:-37:LEU:CD1	2.93	0.51
1:C:57:TYR:CD1	1:C:71:ALA:HB1	2.45	0.51
1:B:23:ARG:HG3	1:B:23:ARG:O	2.11	0.51
1:A:-345:GLU:H	1:A:-342:LYS:HE2	1.75	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:-119:PRO:HA	1:B:-116:TRP:CE2	2.47	0.50
1:B:-338:TRP:CG	1:B:-291:PRO:HG3	2.46	0.50
1:B:-286:TRP:CD1	1:B:-282:ARG:HG3	2.46	0.50
1:A:-195:GLU:OE1	1:A:-4:ARG:NH2	2.44	0.50
1:A:31:GLU:OE1	1:A:35:ARG:CZ	2.60	0.50
1:A:28:ARG:HG3	1:A:28:ARG:HH11	1.76	0.50
1:A:-214:ALA:O	1:A:-210:GLU:HG3	2.11	0.50
1:D:-260:LYS:O	1:D:-259:LEU:HD23	2.12	0.50
1:D:74:VAL:HA	1:D:78:TRP:O	2.12	0.50
1:C:-206:LYS:O	1:C:-206:LYS:HG2	2.11	0.50
1:D:-241:PRO:O	1:D:-63:LEU:HD21	2.11	0.50
1:C:-200:MET:CE	1:C:-126:THR:HB	2.42	0.49
1:D:-227:LEU:CD2	1:D:-204:LYS:HG3	2.43	0.49
1:D:-55:VAL:HG12	1:D:-49:LEU:HG	1.92	0.49
1:D:-130:ASN:OD1	1:D:-113:ILE:HG12	2.12	0.49
1:B:-345:GLU:CD	1:B:-342:LYS:HE2	2.38	0.49
2:E:4:DC:H2''	2:E:5:DT:C6	2.47	0.49
1:C:-219:TRP:O	1:C:-216:ILE:HB	2.12	0.49
2:H:1:DC:H5'	5:H:104:HOH:O	2.11	0.49
1:C:-320:GLU:OE1	1:C:-314:LYS:HD2	2.12	0.49
1:A:36:ARG:HH22	2:F:-1:DC:H3'	1.77	0.49
1:A:-212:ASP:HA	1:A:-202:ALA:HB2	1.94	0.49
1:C:-215:PRO:HA	1:C:-145:HIS:HD2	1.72	0.49
1:D:-346:ILE:HD13	1:D:-77:PRO:CD	2.39	0.49
1:A:31:GLU:HA	1:A:31:GLU:OE2	2.12	0.49
1:B:-226:LEU:HD11	1:B:-213:LEU:HD11	1.95	0.48
1:D:-242:TYR:HD2	1:D:-68:LEU:HD23	1.78	0.48
1:A:30:ARG:HB3	1:A:34:ARG:HH12	1.79	0.48
1:D:-320:GLU:CG	1:D:-314:LYS:HA	2.43	0.48
1:D:-58:LEU:HD23	1:D:-58:LEU:N	2.29	0.48
1:C:-164:ASP:HB3	1:C:17:GLN:OE1	2.14	0.48
1:D:-197:LEU:HD11	1:D:-144:MET:HE3	1.96	0.48
1:A:23:ARG:O	1:A:23:ARG:HG3	2.13	0.48
1:C:-343:GLY:O	1:C:-315:ILE:HG23	2.13	0.47
1:C:-220:THR:HG22	1:C:-99:THR:OG1	2.14	0.47
1:D:-55:VAL:CG1	1:D:-49:LEU:HD21	2.45	0.47
1:D:-315:ILE:CD1	1:D:-73:LEU:CD1	2.73	0.47
1:C:88:ARG:NH2	1:D:75:GLU:OE2	2.44	0.47
1:D:-128:GLY:HA2	5:D:205:HOH:O	2.14	0.47
1:D:-291:PRO:HG2	1:D:-288:ILE:HG23	1.96	0.47
1:D:-276:GLN:HG2	1:D:-276:GLN:O	2.14	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:52:ARG:NH1	1:A:63:CYS:O	2.43	0.47
1:D:-255:THR:HB	1:D:-241:PRO:HB3	1.96	0.47
1:D:-247:GLY:C	1:D:-246:LYS:HD2	2.40	0.47
1:C:-57:GLU:HG3	1:C:-41:TYR:OH	2.16	0.46
1:D:-258:TYR:CE1	1:D:-43:LYS:HA	2.50	0.46
1:D:-19:ILE:HD12	1:D:-19:ILE:C	2.41	0.46
2:F:4:DC:H2''	2:F:5:DT:H5'	1.98	0.46
1:C:-167:VAL:O	1:C:17:GLN:NE2	2.48	0.46
1:D:-191:THR:HG22	1:D:-187:ILE:HD13	1.95	0.46
1:D:-170:ILE:HG21	1:D:-13:GLN:NE2	2.31	0.46
1:C:1:ASN:HB3	1:C:7:GLN:HB2	1.97	0.45
1:C:-263:PHE:CZ	1:C:-67:GLU:HG2	2.51	0.45
1:C:-209:LEU:HD12	1:C:-204:LYS:CB	2.43	0.45
1:A:-345:GLU:N	1:A:-342:LYS:HE3	2.31	0.45
1:B:-194:PRO:HB3	1:B:-5:VAL:HG12	1.98	0.45
1:C:-281:PHE:HB3	1:C:-244:ILE:CD1	2.45	0.45
1:C:-259:LEU:HD23	1:C:-241:PRO:HG2	1.98	0.45
1:D:-191:THR:O	1:D:-187:ILE:CD1	2.64	0.45
1:A:-153:LEU:HD12	1:A:-144:MET:HE1	1.96	0.45
1:C:-194:PRO:HD3	1:C:-4:ARG:CG	2.46	0.45
1:D:-272:LEU:HD22	1:D:-83:GLY:HA3	1.99	0.45
1:A:-119:PRO:HA	1:A:-116:TRP:CE2	2.52	0.45
1:C:-305:LEU:C	1:C:-305:LEU:HD12	2.42	0.45
1:D:-272:LEU:HD13	1:D:-244:ILE:CG2	2.47	0.45
1:D:-180:ALA:O	1:D:-167:VAL:HA	2.17	0.45
1:B:-315:ILE:HG21	1:B:-73:LEU:HD13	1.99	0.45
1:A:46:LYS:HZ2	1:B:85:THR:HG22	1.81	0.44
1:A:-98:PHE:CE2	1:A:-97:LYS:HD2	2.51	0.44
1:C:-269:ILE:HG23	1:C:-242:TYR:CZ	2.52	0.44
1:B:-208:LYS:HD3	1:B:-208:LYS:HA	1.64	0.44
1:C:71:ALA:O	1:C:74:VAL:HG22	2.18	0.44
1:D:-131:PHE:HA	1:D:-126:THR:HG22	2.00	0.44
2:G:-1:DC:H2''	2:G:0:DA:C8	2.53	0.44
1:C:-346:ILE:HG22	1:C:-290:ASP:CG	2.43	0.44
1:C:60:PRO:HG2	1:C:63:CYS:HB3	1.98	0.44
1:D:-227:LEU:HD22	1:D:-204:LYS:HG3	1.99	0.44
1:B:-260:LYS:C	1:B:-259:LEU:HD12	2.42	0.44
1:B:61:LYS:HE2	1:B:62:HIS:CE1	2.53	0.44
1:C:-113:ILE:O	1:C:-110:SER:OG	2.34	0.44
1:D:-324:GLY:O	1:D:-313:VAL:HG23	2.17	0.44
1:C:-213:LEU:HD23	1:C:-213:LEU:O	2.18	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:-191:THR:O	1:D:-187:ILE:HD13	2.18	0.43
1:C:80:VAL:HG23	1:C:80:VAL:O	2.18	0.43
1:D:-219:TRP:CD1	1:D:-100:PRO:HB2	2.53	0.43
1:D:-167:VAL:HG21	1:D:17:GLN:HG3	2.01	0.43
1:C:-177:TYR:OH	1:C:-174:GLY:HA2	2.18	0.43
1:D:-255:THR:HB	1:D:-241:PRO:CB	2.48	0.43
1:C:-60:GLU:H	1:C:-60:GLU:CD	2.26	0.43
1:C:-116:TRP:HB2	1:C:-50:PRO:HG2	2.00	0.43
1:A:-100:PRO:O	1:A:-93:SER:OG	2.30	0.43
1:C:-239:ALA:HA	1:C:-46:VAL:HA	2.00	0.43
1:C:-60:GLU:OE2	1:C:-60:GLU:N	2.46	0.43
1:A:46:LYS:HG2	1:B:84:GLY:C	2.43	0.43
1:B:83:ASP:HB3	1:B:85:THR:HG23	2.01	0.43
1:C:-251:VAL:O	1:C:-244:ILE:HG13	2.18	0.42
1:B:24:LYS:C	1:B:24:LYS:CD	2.92	0.42
1:C:-102:VAL:HG12	1:C:-101:LEU:O	2.19	0.42
1:C:-134:GLU:O	1:C:-133:ALA:C	2.62	0.42
1:C:-232:ILE:HG22	1:C:-131:PHE:CZ	2.55	0.42
1:C:-208:LYS:NZ	1:C:-145:HIS:HA	2.34	0.42
1:D:-300:PRO:HA	1:D:-273:LEU:CD2	2.49	0.42
1:D:-193:TYR:HB2	3:M:2:GLC:O5	2.19	0.42
1:A:27:TRP:O	1:A:27:TRP:HD1	2.03	0.42
1:B:-337:ILE:O	1:B:-309:HIS:HA	2.19	0.42
1:C:-159:LYS:HD3	1:C:10:ASP:OD1	2.19	0.42
1:C:76:ALA:HB1	1:D:75:GLU:HG2	2.01	0.42
1:A:31:GLU:OE1	1:A:35:ARG:NE	2.53	0.42
1:A:81:GLU:CD	1:A:82:GLU:N	2.75	0.42
1:C:-348:MET:O	1:C:-348:MET:HG3	2.17	0.42
1:C:6:ARG:HG3	1:C:6:ARG:HH11	1.84	0.42
1:D:-262:GLN:HA	1:D:-254:TRP:CE2	2.55	0.42
1:A:38:ARG:HA	1:A:41:ARG:CZ	2.49	0.42
1:D:-346:ILE:H	1:D:-346:ILE:HG13	1.55	0.42
1:D:-243:ALA:HB1	1:D:-84:ALA:C	2.45	0.42
1:A:-338:TRP:CD2	1:A:-291:PRO:HG3	2.55	0.42
1:A:27:TRP:O	1:A:27:TRP:CD1	2.71	0.41
1:D:-305:LEU:HD12	1:D:-305:LEU:C	2.45	0.41
2:H:10:DA:H2''	2:H:11:DA:O5'	2.19	0.41
1:C:-188:LEU:O	1:C:-188:LEU:HD12	2.20	0.41
1:D:-308:PRO:HG2	1:D:-305:LEU:HB3	2.02	0.41
1:A:28:ARG:HH11	1:A:28:ARG:CG	2.34	0.41
1:C:-304:GLU:OE1	1:C:-304:GLU:N	2.48	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:F:4:DC:H2''	2:F:5:DT:C5'	2.51	0.41
2:F:10:DA:H2''	2:F:11:DA:O5'	2.21	0.41
1:C:-209:LEU:CD1	1:C:-204:LYS:HB2	2.45	0.41
1:C:-153:LEU:HD11	1:C:-144:MET:HE1	2.03	0.41
1:C:-134:GLU:O	1:C:-130:ASN:OD1	2.38	0.41
1:D:-291:PRO:CG	1:D:-288:ILE:HG23	2.50	0.41
1:D:-271:ALA:HB3	1:D:-82:ILE:HG22	2.02	0.41
1:A:-178:LYS:HA	1:A:-178:LYS:HD3	1.91	0.41
1:C:-180:ALA:O	1:C:-167:VAL:HA	2.21	0.41
1:C:70:LYS:HG2	1:C:80:VAL:HG21	2.02	0.41
1:D:-300:PRO:HA	1:D:-273:LEU:HD21	2.02	0.41
1:D:-251:VAL:HG21	1:D:-243:ALA:HB3	2.03	0.41
1:A:-116:TRP:HB2	1:A:-50:PRO:HG2	2.02	0.41
1:D:-262:GLN:HA	1:D:-254:TRP:NE1	2.36	0.41
1:D:-259:LEU:HB2	1:D:-254:TRP:NE1	2.36	0.41
1:D:-272:LEU:HD13	1:D:-244:ILE:HG21	2.03	0.41
1:D:-254:TRP:CE3	1:D:-245:LEU:CD2	3.04	0.40
1:C:-319:LYS:HE3	1:C:-319:LYS:HB3	1.76	0.40
1:C:-206:LYS:O	1:C:-206:LYS:CG	2.70	0.40
1:D:-233:LEU:N	1:D:-101:LEU:HD23	2.36	0.40
1:D:-51:LYS:HD2	1:D:-51:LYS:N	2.35	0.40
1:C:-184:ASP:HB3	1:C:-161:GLY:HA2	2.02	0.40
1:A:-338:TRP:CG	1:A:-291:PRO:HG3	2.56	0.40
1:D:-220:THR:OG1	1:D:-218:GLU:OE1	2.39	0.40
1:B:-131:PHE:HA	1:B:-126:THR:HG22	2.04	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	435/453 (96%)	429 (99%)	6 (1%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	B	434/453 (96%)	429 (99%)	5 (1%)	0	100	100
1	C	435/453 (96%)	425 (98%)	10 (2%)	0	100	100
1	D	434/453 (96%)	424 (98%)	10 (2%)	0	100	100
All	All	1738/1812 (96%)	1707 (98%)	31 (2%)	0	100	100

There are no Ramachandran outliers to report.

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	344/356 (97%)	333 (97%)	11 (3%)	34	40
1	B	343/356 (96%)	336 (98%)	7 (2%)	48	58
1	C	344/356 (97%)	333 (97%)	11 (3%)	34	40
1	D	343/356 (96%)	327 (95%)	16 (5%)	23	24
All	All	1374/1424 (96%)	1329 (97%)	45 (3%)	33	39

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

Mol	Chain	Res	Type
1	C	-346	ILE
1	C	-295	THR
1	C	-227	LEU
1	C	-216	ILE
1	C	-213	LEU
1	C	-209	LEU
1	C	-146	LYS
1	C	-143	ASN
1	C	-90	PHE
1	C	-42	SER
1	C	17	GLN
1	D	-346	ILE
1	D	-337	ILE

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Mol	Chain	Res	Type
1	D	-317	THR
1	D	-315	ILE
1	D	-314	LYS
1	D	-313	VAL
1	D	-311	VAL
1	D	-269	ILE
1	D	-111	THR
1	D	-103	THR
1	D	-90	PHE
1	D	-82	ILE
1	D	-73	LEU
1	D	-42	SER
1	D	23	ARG
1	D	88	ARG
1	A	-347	LYS
1	A	-346	ILE
1	A	-345	GLU
1	A	-342	LYS
1	A	-288	ILE
1	A	-90	PHE
1	A	17	GLN
1	A	41	ARG
1	A	46	LYS
1	A	79	VAL
1	A	80	VAL
1	B	-347	LYS
1	B	-115	SER
1	B	-90	PHE
1	B	28	ARG
1	B	58	ASN
1	B	85	THR
1	B	88	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (11) such sidechains are listed below:

Mol	Chain	Res	Type
1	C	-224	ASN
1	C	-54	ASN
1	C	54	GLN
1	D	-276	GLN
1	D	-262	GLN
1	D	62	HIS

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Mol	Chain	Res	Type
1	A	-336	ASN
1	B	-336	ASN
1	B	-330	ASN
1	B	-276	GLN
1	B	-224	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 ⓘ

8 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	$\# Z > 2$
3	GLC	J	1	3	12,12,12	0.48	0	17,17,17	0.50	0
3	GLC	J	2	3	11,11,12	0.55	0	15,15,17	0.88	0
3	GLC	K	1	3	12,12,12	0.50	0	17,17,17	0.72	0
3	GLC	K	2	3	11,11,12	0.57	0	15,15,17	0.72	0
3	GLC	L	1	3	12,12,12	0.49	0	17,17,17	0.63	0
3	GLC	L	2	3	11,11,12	0.53	0	15,15,17	0.77	0
3	GLC	M	1	3	12,12,12	0.53	0	17,17,17	0.58	0
3	GLC	M	2	3	11,11,12	0.59	0	15,15,17	0.62	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	GLC	J	1	3	-	0/2/22/22	0/1/1/1
3	GLC	J	2	3	-	0/2/19/22	0/1/1/1
3	GLC	K	1	3	-	0/2/22/22	0/1/1/1
3	GLC	K	2	3	-	0/2/19/22	0/1/1/1
3	GLC	L	1	3	-	0/2/22/22	0/1/1/1
3	GLC	L	2	3	-	0/2/19/22	0/1/1/1
3	GLC	M	1	3	-	2/2/22/22	0/1/1/1
3	GLC	M	2	3	-	0/2/19/22	0/1/1/1

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (2) torsion outliers are listed below:

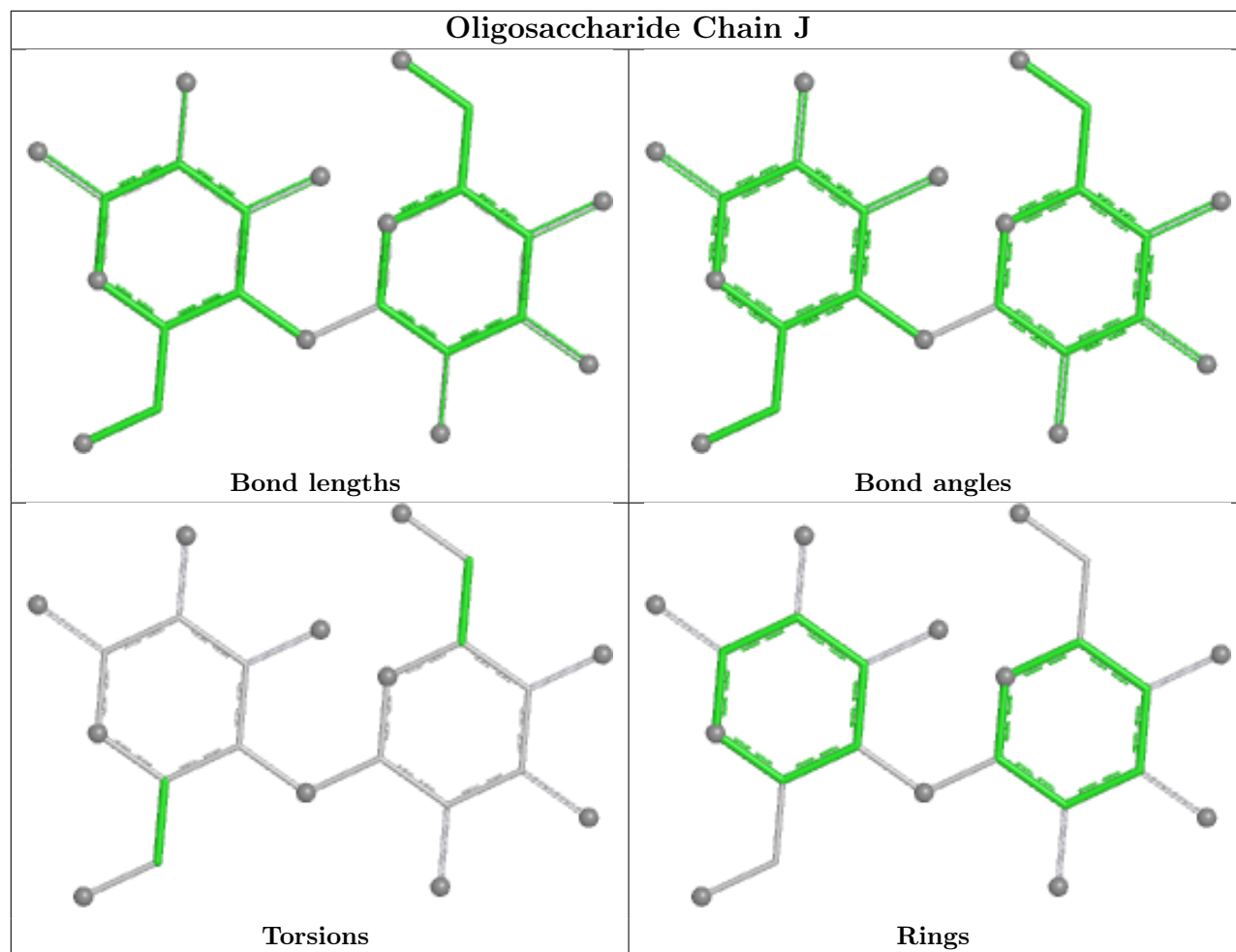
Mol	Chain	Res	Type	Atoms
3	M	1	GLC	C4-C5-C6-O6
3	M	1	GLC	O5-C5-C6-O6

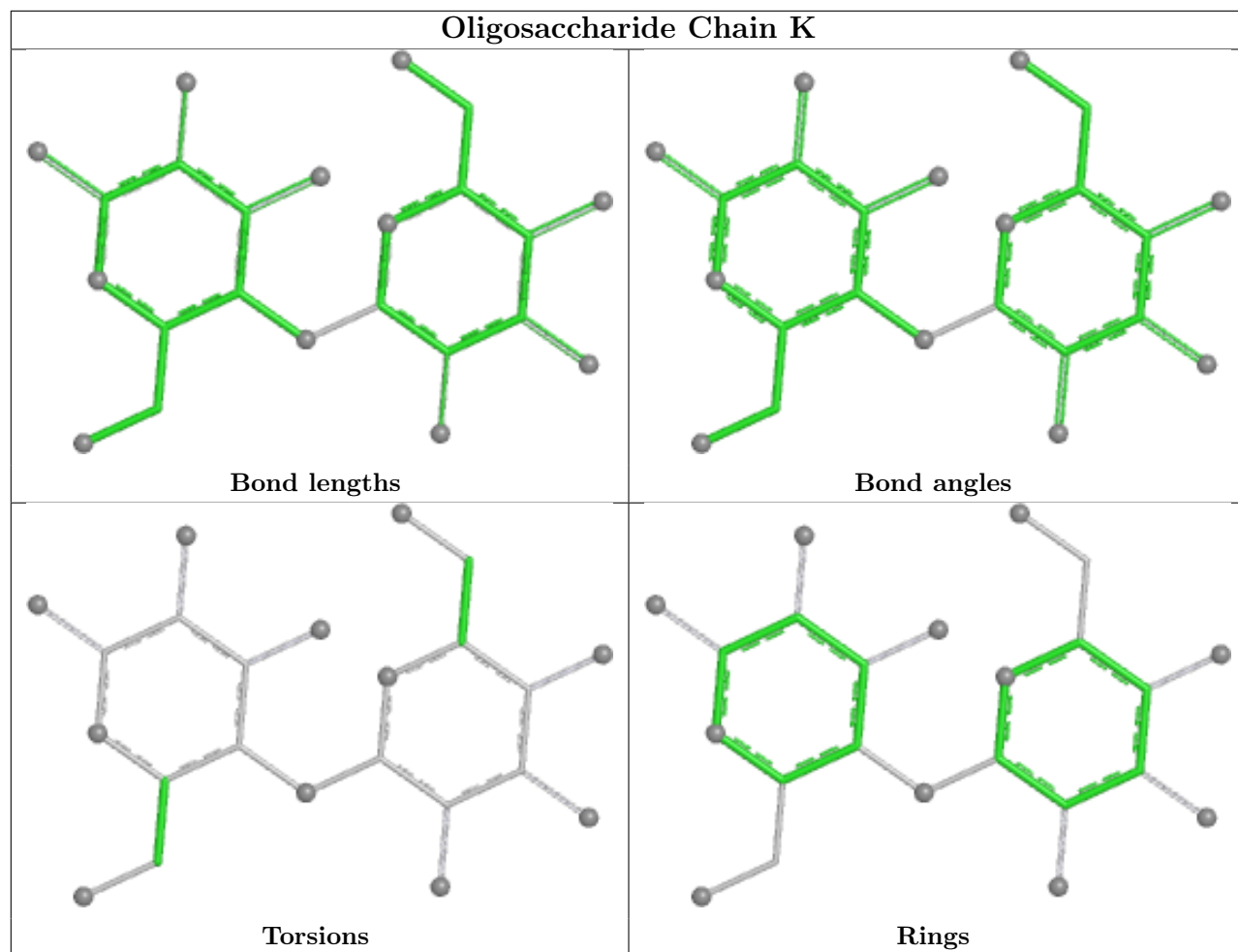
There are no ring outliers.

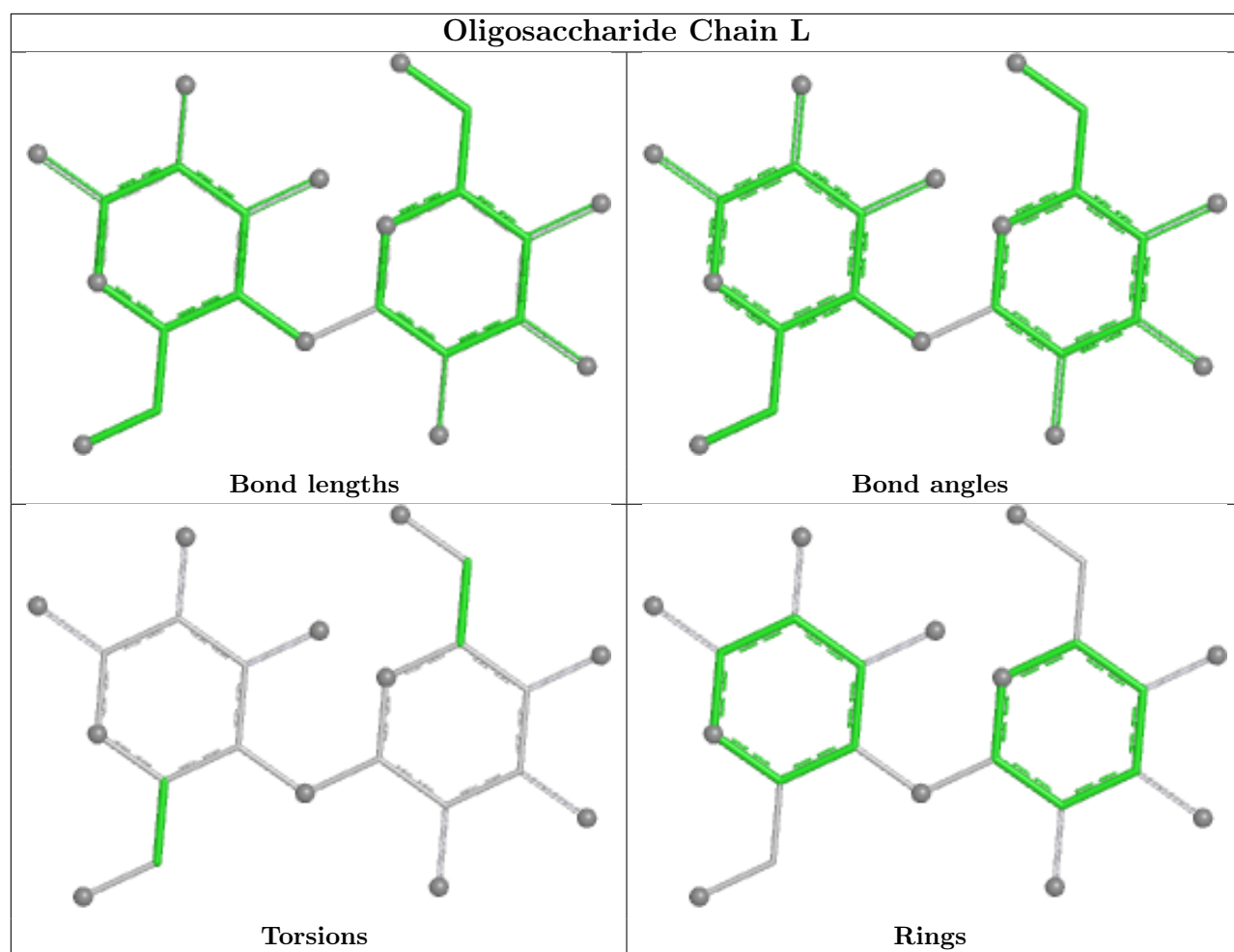
1 monomer is involved in 1 short contact:

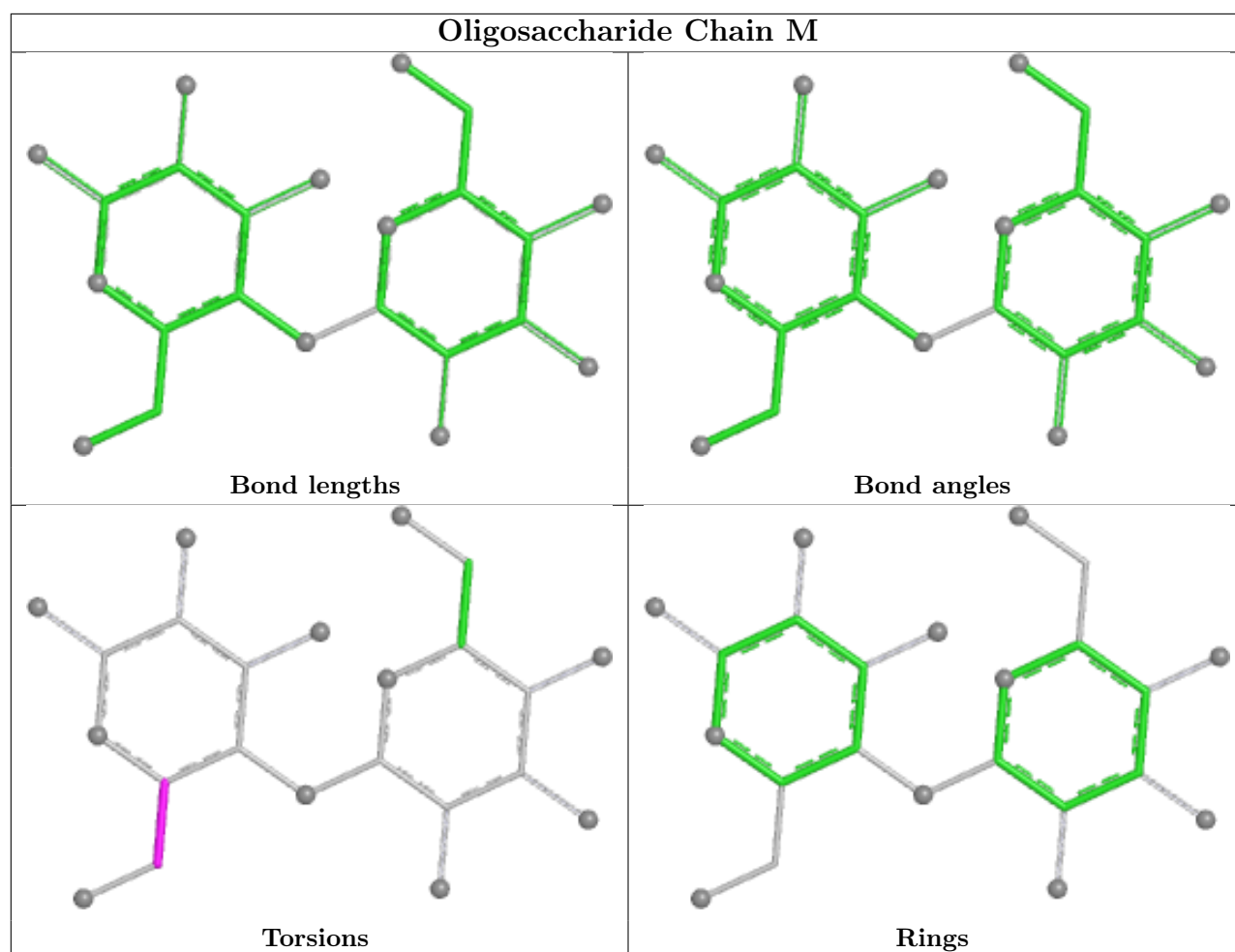
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	M	2	GLC	1	0

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









5.6 Ligand geometry [i](#)

3 ligands 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	$\# Z > 2$
4	EDO	A	301	-	3,3,3	0.43	0	2,2,2	0.45	0
4	EDO	C	301	-	3,3,3	0.44	0	2,2,2	0.61	0
4	EDO	B	301	-	3,3,3	0.43	0	2,2,2	0.43	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the

Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns.
'-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	EDO	A	301	-	-	1/1/1/1	-
4	EDO	C	301	-	-	0/1/1/1	-
4	EDO	B	301	-	-	1/1/1/1	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	B	301	EDO	O1-C1-C2-O2
4	A	301	EDO	O1-C1-C2-O2

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 ⓘ

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	437/453 (96%)	0.16	14 (3%) 50 50	26, 45, 76, 110	0
1	B	436/453 (96%)	0.04	11 (2%) 58 59	22, 40, 74, 83	0
1	C	437/453 (96%)	0.31	9 (2%) 63 65	26, 53, 88, 127	0
1	D	436/453 (96%)	0.81	54 (12%) 8 7	36, 65, 104, 119	0
2	E	15/15 (100%)	0.29	0 100 100	55, 64, 79, 85	0
2	F	15/15 (100%)	0.67	1 (6%) 24 22	40, 75, 95, 105	0
2	G	15/15 (100%)	0.19	0 100 100	42, 55, 75, 78	0
2	H	15/15 (100%)	-0.14	0 100 100	47, 52, 60, 62	0
All	All	1806/1872 (96%)	0.33	89 (4%) 35 33	22, 52, 91, 127	0

All (89) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	27	TRP	5.2
1	D	-266	ALA	4.2
1	A	-7	TYR	3.9
1	B	-27	MET	3.7
1	D	-346	ILE	3.6
1	D	-73	LEU	3.6
1	D	-287	PHE	3.5
1	D	-317	THR	3.5
1	C	-109	ALA	3.5
1	C	-207	ALA	3.4
1	D	-82	ILE	3.3
1	A	36	ARG	3.2
1	D	-288	ILE	3.2
1	A	33	ASN	3.0
1	D	-244	ILE	3.0
1	D	-267	PRO	3.0

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Mol	Chain	Res	Type	RSRZ
1	D	-68	LEU	2.9
1	D	-7	TYR	2.8
1	A	42	ALA	2.8
1	C	-214	ALA	2.8
1	B	-7	TYR	2.8
1	D	-265	ALA	2.8
1	D	-341	LEU	2.7
1	B	27	TRP	2.7
1	A	28	ARG	2.7
1	C	-108	VAL	2.6
1	D	-176	ALA	2.6
1	B	32	ASN	2.6
1	D	-311	VAL	2.6
2	F	-3	DT	2.5
1	D	-314	LYS	2.5
1	D	-271	ALA	2.5
1	D	-343	GLY	2.5
1	B	35	ARG	2.4
1	C	-128	GLY	2.4
1	D	-258	TYR	2.4
1	D	-242	TYR	2.4
1	A	57	TYR	2.3
1	D	-316	GLY	2.3
1	A	23	ARG	2.3
1	B	24	LYS	2.3
1	D	26	SER	2.3
1	D	-260	LYS	2.3
1	D	-257	PRO	2.3
1	D	-252	ALA	2.3
1	D	-345	GLU	2.3
1	D	-289	ILE	2.3
1	D	-313	VAL	2.3
1	C	-348	MET	2.2
1	B	-294	GLY	2.2
1	D	-327	ALA	2.2
1	A	-176	ALA	2.2
1	D	-240	ILE	2.2
1	D	-296	ALA	2.2
1	A	24	LYS	2.2
1	D	-256	PHE	2.2
1	D	-248	ASN	2.2
1	D	-312	THR	2.2

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Mol	Chain	Res	Type	RSRZ
1	D	-79	ALA	2.2
1	D	-81	ASN	2.2
1	D	-249	TYR	2.1
1	D	-328	LEU	2.1
1	D	-259	LEU	2.1
1	D	-63	LEU	2.1
1	D	-141	ASP	2.1
1	B	64	ASP	2.1
1	D	-342	LYS	2.1
1	A	29	GLU	2.1
1	D	-277	ALA	2.1
1	C	-227	LEU	2.1
1	D	-273	LEU	2.1
1	A	56	ASP	2.1
1	D	-329	GLY	2.1
1	B	23	ARG	2.1
1	C	-7	TYR	2.1
1	D	-261	ASP	2.1
1	B	60	PRO	2.1
1	A	26	SER	2.1
1	B	-296	ALA	2.1
1	D	-301	PHE	2.1
1	A	41	ARG	2.1
1	D	-84	ALA	2.0
1	D	-315	ILE	2.0
1	D	-41	TYR	2.0
1	D	-322	LYS	2.0
1	C	-265	ALA	2.0
1	D	-264	ALA	2.0
1	D	-66	ASN	2.0
1	D	-269	ILE	2.0

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

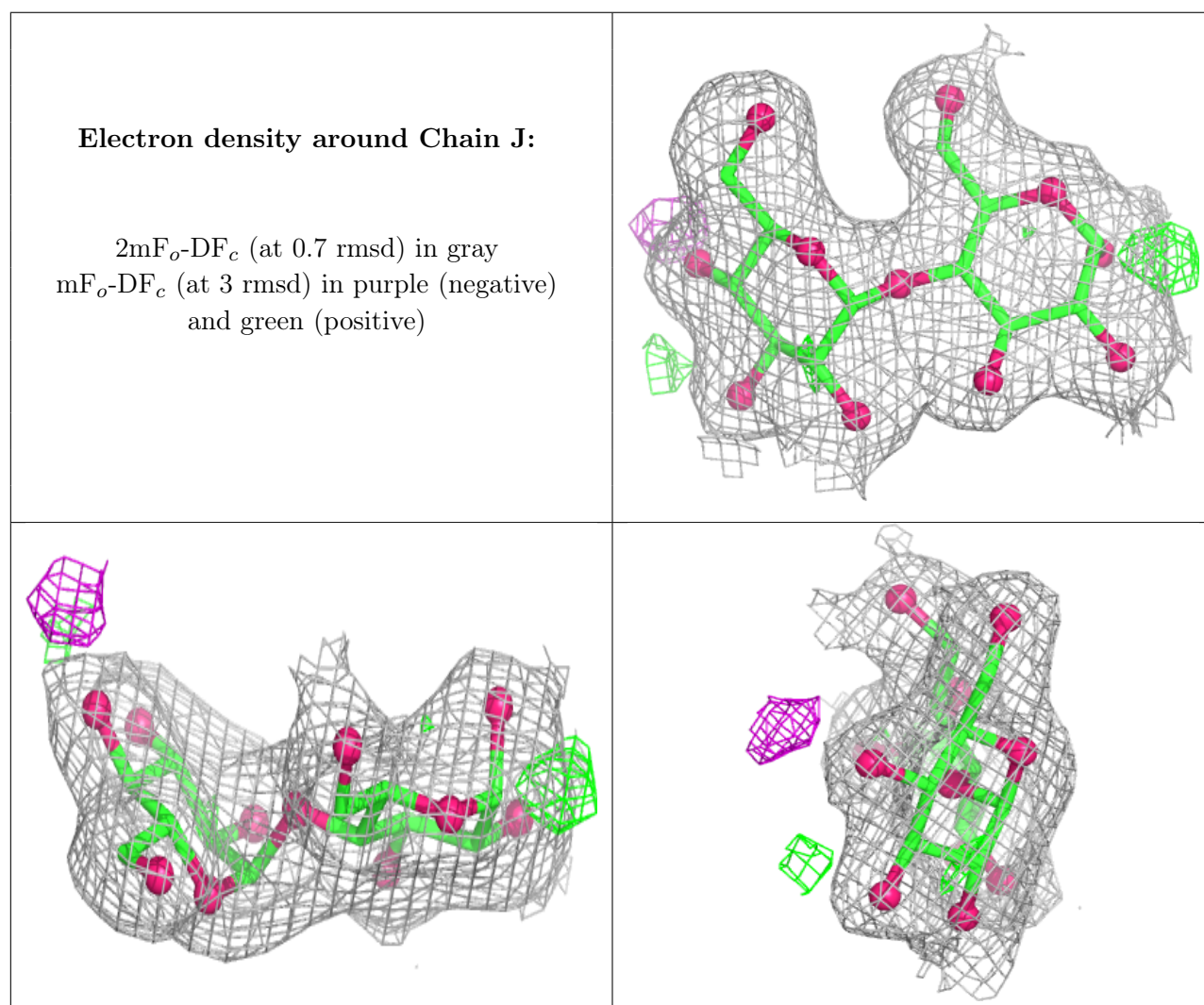
There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates [i](#)

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

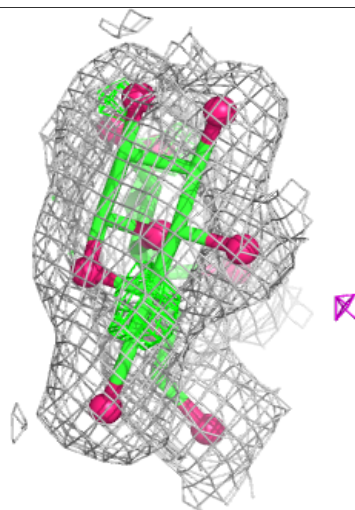
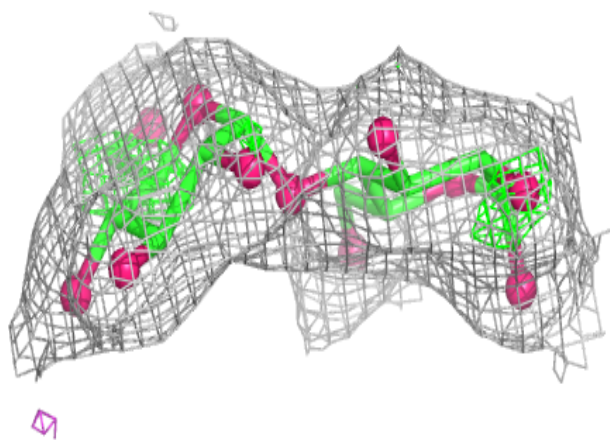
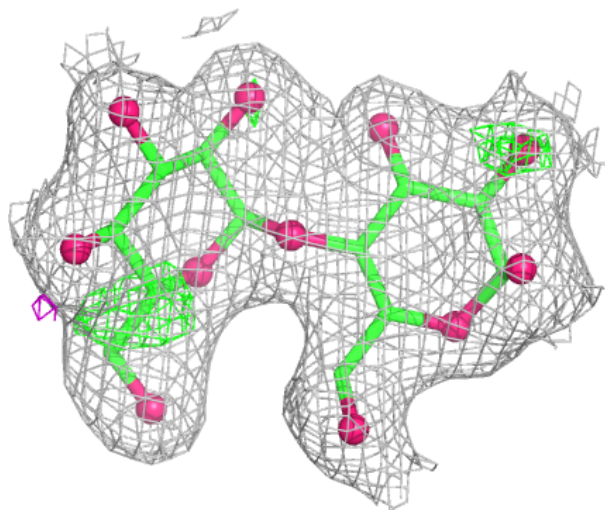
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	GLC	J	1	12/12	0.93	0.09	29,33,36,45	0
3	GLC	L	2	11/12	0.93	0.09	26,36,41,46	0
3	GLC	K	1	12/12	0.95	0.07	23,32,37,50	0
3	GLC	L	1	12/12	0.95	0.08	29,38,42,48	0
3	GLC	J	2	11/12	0.95	0.08	25,31,33,38	0
3	GLC	K	2	11/12	0.96	0.07	20,29,41,43	0
3	GLC	M	1	12/12	-	-	34,48,59,65	0
3	GLC	M	2	11/12	-	-	44,47,60,62	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.



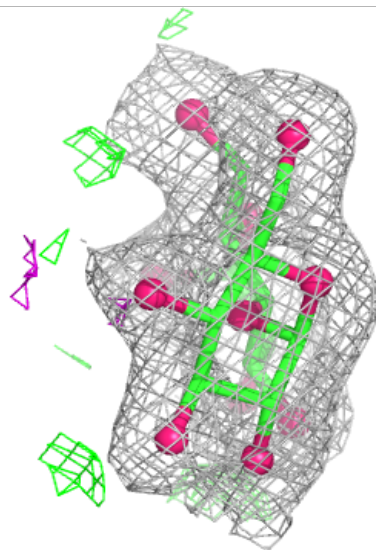
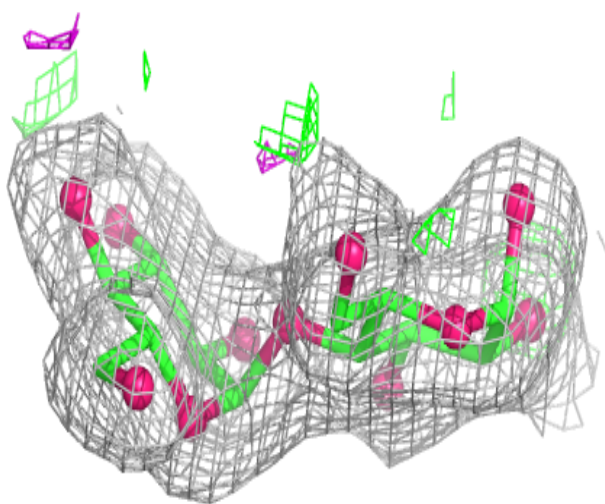
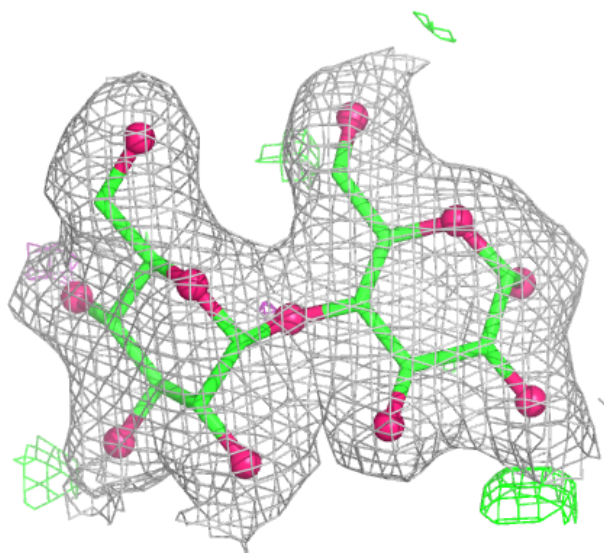
Electron density around Chain K:

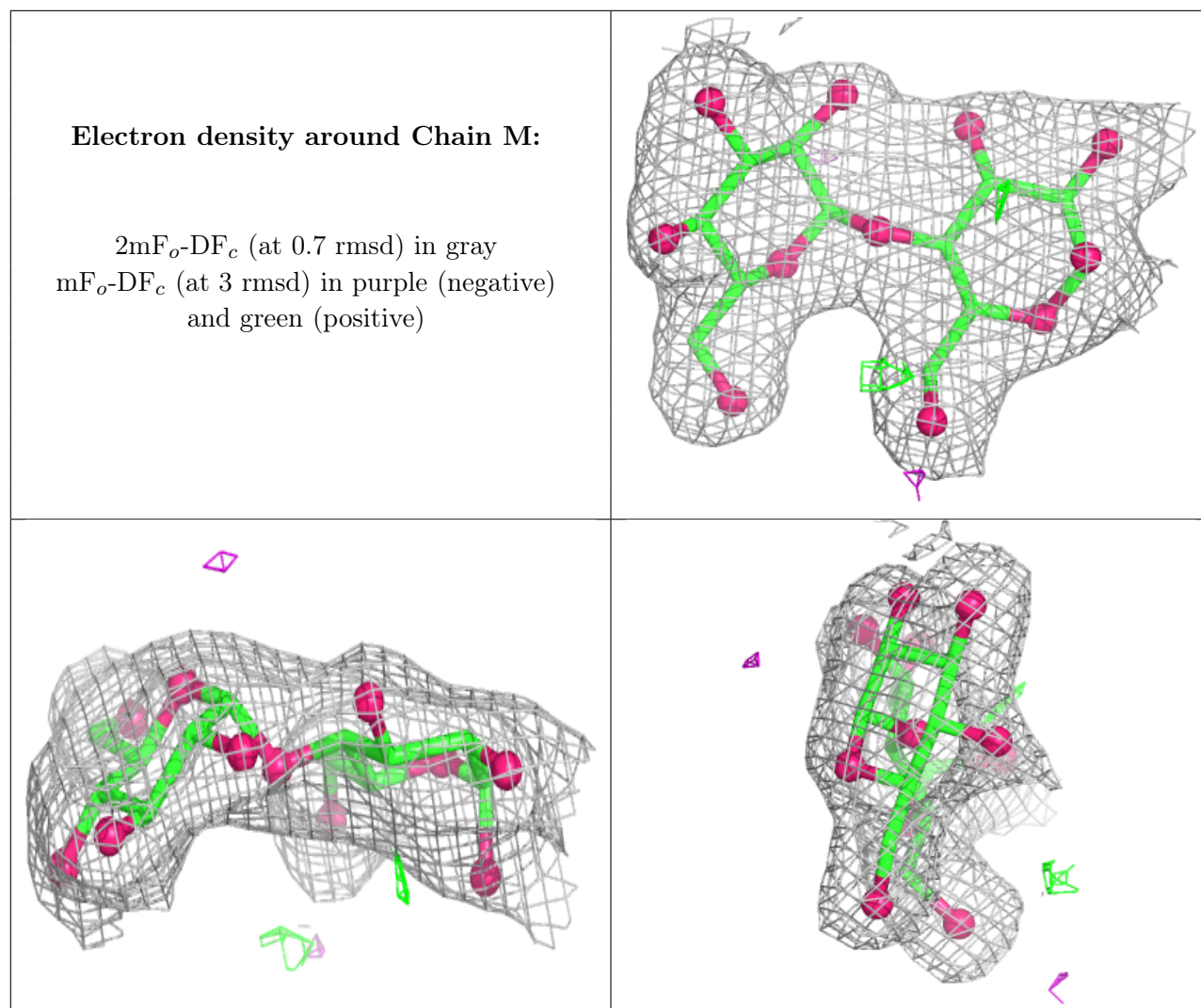
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



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	EDO	B	301	4/4	0.90	0.15	37,45,53,53	0
4	EDO	C	301	4/4	0.94	0.10	42,51,58,58	0
4	EDO	A	301	4/4	0.95	0.10	37,45,52,52	0

6.5 Other polymers [i](#)

There are no such residues in this entry.