



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 20, 2026 – 11:38 AM UTC

PDB ID : 1CXF / pdb_00001cxf
Title : COMPLEX OF A (D229N/E257Q) DOUBLE MUTANT CGTASE FROM BACILLUS CIRCULANS STRAIN 251 WITH MALTOTETRAOSE AT 120 K AND PH 9.1 OBTAINED AFTER SOAKING THE CRYSTAL WITH ALPHA-CYCLODEXTRIN
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Deposited on : 1995-07-28
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

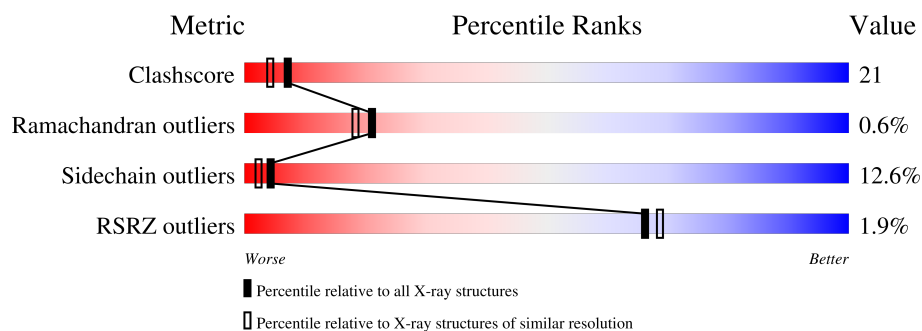
1 Overall quality at a glance

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))
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	686	
2	B	4	
3	C	2	
4	D	6	
4	E	6	

2 Entry composition [i](#)

There are 6 unique types of molecules in this entry. The entry contains 5776 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 CYCLODEXTRIN GLYCOSYLTRANSFERASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	686	5264	3321	902	1025	16	0	0	0

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	229	ASN	ASP	conflict	UNP P43379
A	257	GLN	GLU	conflict	UNP P43379

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



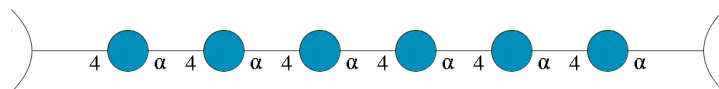
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace
			Total	C	O			
2	B	4	45	24	21	0	0	0

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



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

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



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace
4	D	6	Total	C	O	0	0	0
			66	36	30			
4	E	6	Total	C	O	0	0	0
			66	36	30			

- Molecule 5 is CALCIUM ION (CCD ID: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	2	Total	Ca	0	0
			2	2		

- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	310	Total	O	0	0
			310	310		

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

Chain C:  100%

GLC1
GLC2

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

Chain D:  100%

GLC1
GLC2
GLC3
GLC4
GLC5
GLC6

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

Chain E:  50% 33% 17%

GLC1
GLC2
GLC3
GLC4
GLC5
GLC6

4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	116.84Å 110.00Å 67.42Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	8.00 – 2.10 8.00 – 2.62	Depositor EDS
% Data completeness (in resolution range)	(Not available) (8.00-2.10) 20.0 (8.00-2.62)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.37 (at 2.61Å)	Xtriage
Refinement program	TNT	Depositor
R, R_{free}	0.190 , 0.200 0.170 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å ²)	21.8	Xtriage
Anisotropy	2.766	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 67.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	5776	wwPDB-VP
Average B, all atoms (Å ²)	11.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.39% 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: CA, 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	0.74	0/5394	1.05	23/7352 (0.3%)

There are no bond length outliers.

All (23) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	609	SER	N-CA-C	8.29	122.82	112.87
1	A	546	ASP	N-CA-C	7.71	120.66	111.33
1	A	216	LYS	N-CA-C	-7.25	103.45	111.36
1	A	164	LEU	N-CA-C	-6.97	102.76	112.45
1	A	475	GLY	N-CA-C	-6.94	104.39	115.08
1	A	330	GLU	N-CA-C	-6.52	101.48	110.35
1	A	407	GLY	N-CA-C	6.28	120.67	112.25
1	A	103	ARG	N-CA-C	-6.15	107.61	114.62
1	A	663	GLU	N-CA-C	-6.12	102.41	110.43
1	A	383	THR	N-CA-C	-6.10	105.70	113.02
1	A	473	SER	N-CA-C	-5.82	103.25	110.41
1	A	20	ILE	N-CA-C	5.82	116.24	107.80
1	A	656	GLN	N-CA-C	-5.81	97.73	107.32
1	A	660	VAL	N-CA-C	5.62	117.50	108.85
1	A	167	TYR	N-CA-C	5.58	117.36	111.28
1	A	532	GLY	N-CA-C	-5.41	104.85	111.35
1	A	44	THR	N-CA-C	-5.38	106.88	113.50
1	A	642	VAL	N-CA-C	5.37	113.80	107.84
1	A	460	LEU	N-CA-C	-5.29	106.85	113.15
1	A	53	ASP	N-CA-C	5.17	115.88	108.74
1	A	333	HIS	N-CA-C	5.08	117.04	108.96
1	A	174	LEU	N-CA-C	-5.08	106.60	112.89
1	A	64	GLY	N-CA-C	5.02	122.03	115.40

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	5264	0	5029	219	0
2	B	45	0	39	5	0
3	C	23	0	21	0	0
4	D	66	0	54	0	0
4	E	66	0	54	1	0
5	A	2	0	0	0	0
6	A	310	0	0	7	0
All	All	5776	0	5197	222	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 21.

All (222) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:643:PRO:HB2	1:A:646:LYS:HG3	1.23	1.16
1:A:300:MET:HG3	1:A:419:LEU:HB2	1.25	1.09
1:A:358:ILE:HD13	1:A:391:ILE:HD13	1.35	1.08
1:A:617:ASP:HB3	1:A:620:LYS:HE2	1.47	0.94
1:A:339:ARG:HG2	1:A:365:TYR:CD2	2.04	0.92
1:A:136:PHE:CE2	1:A:138:PRO:HG3	2.06	0.91
1:A:643:PRO:CB	1:A:646:LYS:HG3	2.02	0.88
1:A:12:PHE:CE2	1:A:133:ILE:HD11	2.08	0.88
1:A:617:ASP:CB	1:A:620:LYS:HE2	2.09	0.83
1:A:673:PRO:HG3	1:A:678:ALA:HB2	1.60	0.82
1:A:562:ASN:HD22	1:A:575:VAL:HG11	1.44	0.81
1:A:147:ASP:O	1:A:149:PRO:HD3	1.82	0.80
1:A:358:ILE:CD1	1:A:391:ILE:HD13	2.13	0.79
1:A:235:PRO:O	1:A:239:GLN:HG3	1.83	0.78
1:A:187:GLU:CD	1:A:628:GLN:HB2	2.09	0.77
1:A:333:HIS:O	1:A:367:SER:HB3	1.86	0.76

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:136:PHE:CD2	1:A:138:PRO:HG3	2.22	0.74
1:A:54:TRP:O	1:A:58:ILE:HG13	1.87	0.74
1:A:26:SER:O	1:A:56:GLY:HA3	1.87	0.73
1:A:115:ILE:O	1:A:119:GLN:HG3	1.87	0.73
1:A:600:LEU:HD13	4:E:4:GLC:H3	1.70	0.73
1:A:186:THR:O	1:A:190:ILE:HG13	1.89	0.72
1:A:88:ASN:HD21	1:A:91:GLY:H	1.40	0.70
1:A:163:LEU:HD22	1:A:165:GLY:N	2.06	0.69
1:A:227:ARG:NH1	1:A:257:GLN:HB2	2.08	0.69
1:A:241:SER:HB3	1:A:641:SER:HB2	1.75	0.69
1:A:81:GLU:OE1	1:A:103:ARG:HD3	1.93	0.68
1:A:519:GLY:O	1:A:520:ARG:HD3	1.94	0.67
1:A:300:MET:HE2	1:A:417:ASP:O	1.94	0.67
1:A:339:ARG:HG2	1:A:365:TYR:CE2	2.30	0.66
2:B:2:GLC:H61	2:B:3:GLC:H3	1.77	0.66
1:A:370:THR:HB	6:A:848:HOH:O	1.95	0.66
1:A:424:LYS:HG3	1:A:429:VAL:HG22	1.78	0.66
1:A:598:THR:HB	1:A:602:GLN:HG2	1.77	0.65
1:A:271:LYS:HG3	1:A:275:GLU:OE2	1.95	0.65
1:A:21:PHE:HB2	1:A:360:TYR:CE1	2.31	0.65
1:A:544:TRP:CZ3	1:A:549:ILE:HD13	2.31	0.64
1:A:645:GLY:HA2	1:A:672:ALA:O	1.97	0.64
1:A:332:PHE:CE1	1:A:342:LEU:HA	2.34	0.63
1:A:81:GLU:CD	1:A:103:ARG:HD3	2.22	0.63
1:A:116:ALA:HA	1:A:119:GLN:HE21	1.64	0.62
1:A:434:VAL:HG22	1:A:487:THR:HG23	1.81	0.62
1:A:249:TYR:CE2	1:A:250:LYS:HD2	2.35	0.62
1:A:655:LYS:HE3	6:A:808:HOH:O	1.99	0.62
1:A:214:ALA:O	1:A:217:MET:HB3	2.00	0.62
1:A:17:ILE:HD12	1:A:355:VAL:HG12	1.80	0.61
1:A:24:ARG:HD2	1:A:375:ARG:O	2.00	0.61
1:A:300:MET:CG	1:A:419:LEU:HB2	2.17	0.61
1:A:232:LYS:HD2	1:A:258:TRP:CE2	2.36	0.61
1:A:163:LEU:HD22	1:A:165:GLY:H	1.65	0.61
1:A:142:SER:OG	1:A:155:GLY:HA2	2.01	0.60
1:A:80:VAL:HB	1:A:105:PHE:HA	1.82	0.60
1:A:560:ASN:HD21	1:A:578:ASN:HA	1.67	0.60
1:A:227:ARG:C	1:A:227:ARG:HD2	2.26	0.59
1:A:300:MET:HG3	1:A:419:LEU:CB	2.17	0.59
1:A:14:THR:O	1:A:399:LYS:HG2	2.02	0.59
1:A:607:THR:HG21	1:A:653:LEU:HD21	1.85	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:290:ARG:HH21	1:A:325:ASP:CB	2.16	0.58
1:A:643:PRO:HB2	1:A:646:LYS:CG	2.16	0.58
1:A:70:GLY:O	1:A:72:THR:HG23	2.03	0.58
1:A:101:TRP:HE1	2:B:4:GLC:H61	1.68	0.58
1:A:361:GLY:HA3	1:A:366:MET:SD	2.44	0.58
1:A:395:ALA:HB3	1:A:396:PRO:HD3	1.86	0.58
1:A:633:TYR:CG	1:A:634:PRO:HA	2.38	0.58
1:A:326:ASN:OD1	1:A:329:MET:HG3	2.05	0.57
1:A:583:SER:HB2	1:A:643:PRO:HG3	1.85	0.57
1:A:291:GLN:HG2	1:A:297:THR:OG1	2.04	0.56
1:A:401:ASN:OD1	1:A:403:ALA:HB3	2.05	0.56
1:A:453:GLN:HA	1:A:471:VAL:O	2.05	0.56
1:A:590:ARG:HG2	1:A:679:THR:HG23	1.87	0.56
1:A:630:VAL:HG12	1:A:631:TYR:CD1	2.41	0.56
1:A:290:ARG:HH21	1:A:325:ASP:HB3	1.71	0.56
1:A:518:ASP:OD1	1:A:548:GLN:HG3	2.05	0.56
1:A:273:ALA:HB2	1:A:280:LEU:HD12	1.87	0.55
1:A:562:ASN:ND2	1:A:575:VAL:HG11	2.17	0.55
1:A:649:GLU:HA	1:A:668:HIS:O	2.06	0.55
1:A:3:ASP:HA	1:A:548:GLN:NE2	2.22	0.54
1:A:12:PHE:O	1:A:355:VAL:HG13	2.07	0.54
1:A:387:ALA:O	1:A:391:ILE:HG13	2.07	0.54
1:A:187:GLU:OE1	1:A:628:GLN:HB2	2.07	0.54
1:A:34:PRO:HG2	1:A:49:TYR:CG	2.43	0.54
1:A:40:ASP:OD1	1:A:42:THR:HG22	2.08	0.54
1:A:227:ARG:HD2	1:A:227:ARG:O	2.08	0.54
1:A:187:GLU:HB2	1:A:626:TYR:CD1	2.43	0.54
1:A:187:GLU:OE2	1:A:628:GLN:HB2	2.07	0.54
1:A:204:ASN:OD1	1:A:206:THR:HB	2.07	0.53
1:A:648:ILE:HB	1:A:670:PHE:CE1	2.43	0.53
1:A:655:LYS:HG2	1:A:660:VAL:HB	1.92	0.52
1:A:358:ILE:HD13	1:A:391:ILE:CD1	2.23	0.52
1:A:341:LYS:HD2	6:A:898:HOH:O	2.09	0.51
1:A:83:ILE:HD13	1:A:142:SER:HB2	1.92	0.51
1:A:342:LEU:HD11	1:A:362:THR:HG23	1.92	0.51
1:A:633:TYR:CD2	1:A:634:PRO:HA	2.45	0.51
1:A:271:LYS:O	1:A:275:GLU:HG2	2.10	0.51
1:A:60:LYS:HE2	1:A:381:PHE:CD1	2.45	0.51
1:A:157:LEU:O	1:A:163:LEU:HD23	2.10	0.51
1:A:33:ASN:ND2	1:A:39:PHE:CZ	2.79	0.51
2:B:2:GLC:H61	2:B:3:GLC:H5	1.91	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:627:ASN:HB2	1:A:633:TYR:HB2	1.92	0.51
1:A:655:LYS:HA	1:A:659:THR:O	2.10	0.51
1:A:423:ARG:HH11	1:A:423:ARG:HG3	1.77	0.50
1:A:566:ALA:HA	1:A:571:THR:O	2.11	0.50
1:A:50:CYS:HB3	1:A:377:ARG:NH2	2.27	0.50
1:A:466:GLY:CA	1:A:488:ALA:HB2	2.42	0.50
1:A:561:TYR:CD1	1:A:561:TYR:N	2.79	0.50
1:A:618:PRO:HB3	1:A:653:LEU:HD13	1.94	0.49
1:A:212:LYS:O	1:A:216:LYS:HG2	2.11	0.49
1:A:50:CYS:HB3	1:A:377:ARG:HH21	1.77	0.49
1:A:563:ILE:HD11	1:A:579:PHE:CB	2.42	0.49
1:A:626:TYR:O	1:A:636:TRP:HA	2.13	0.49
1:A:544:TRP:CE3	1:A:549:ILE:HD13	2.47	0.49
1:A:563:ILE:N	1:A:563:ILE:HD12	2.28	0.48
1:A:663:GLU:HA	1:A:684:TRP:CZ2	2.49	0.48
1:A:598:THR:HG22	1:A:602:GLN:CG	2.42	0.48
1:A:378:ILE:O	1:A:378:ILE:HG23	2.12	0.48
1:A:42:THR:CG2	1:A:44:THR:HB	2.44	0.48
1:A:397:LEU:HA	1:A:400:CYS:SG	2.54	0.48
1:A:74:ILE:HD12	1:A:76:ILE:CG2	2.42	0.48
1:A:471:VAL:CG1	1:A:475:GLY:HA2	2.43	0.48
1:A:241:SER:CB	1:A:641:SER:HB2	2.41	0.48
1:A:16:VAL:HG21	1:A:395:ALA:CB	2.43	0.48
1:A:83:ILE:HG22	1:A:153:GLU:OE1	2.14	0.48
1:A:627:ASN:ND2	1:A:636:TRP:CZ2	2.82	0.48
1:A:306:MET:HE3	1:A:307:LEU:HG	1.95	0.48
1:A:340:ARG:O	1:A:340:ARG:HD3	2.13	0.48
1:A:371:ASP:HA	1:A:372:PRO:HA	1.53	0.47
1:A:12:PHE:CE1	1:A:253:PHE:HB3	2.50	0.47
1:A:293:PHE:HE1	1:A:434:VAL:HG11	1.80	0.47
1:A:593:VAL:HA	1:A:682:VAL:O	2.15	0.47
1:A:424:LYS:CG	1:A:429:VAL:HG22	2.45	0.47
1:A:8:ASN:ND2	1:A:11:ASN:HB3	2.30	0.47
1:A:445:SER:HA	1:A:479:ASN:HD22	1.80	0.47
1:A:136:PHE:C	1:A:138:PRO:HD3	2.40	0.46
1:A:505:GLY:HA2	1:A:506:PRO:C	2.41	0.46
1:A:632:GLN:O	1:A:635:ASN:HB2	2.15	0.46
1:A:69:MET:HG3	1:A:388:TYR:CE2	2.51	0.46
1:A:148:GLN:NE2	1:A:148:GLN:HA	2.30	0.46
1:A:290:ARG:NH2	1:A:325:ASP:HB3	2.30	0.46
1:A:563:ILE:N	1:A:563:ILE:CD1	2.78	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:587:VAL:HG21	1:A:673:PRO:HD2	1.97	0.46
1:A:653:LEU:HD21	1:A:662:TRP:CZ3	2.50	0.46
1:A:528:THR:HG22	1:A:529:VAL:N	2.31	0.46
1:A:96:ALA:HB2	1:A:101:TRP:CE3	2.51	0.45
1:A:365:TYR:OH	1:A:386:THR:HB	2.16	0.45
1:A:397:LEU:HD13	1:A:404:ILE:HD13	1.98	0.45
1:A:60:LYS:HE2	1:A:381:PHE:CE1	2.52	0.45
1:A:303:LEU:O	1:A:306:MET:HB3	2.16	0.45
1:A:627:ASN:HB2	1:A:633:TYR:CA	2.46	0.45
1:A:607:THR:HG22	1:A:653:LEU:HG	1.99	0.45
1:A:326:ASN:HB3	6:A:738:HOH:O	2.16	0.45
1:A:374:ASN:OD1	1:A:375:ARG:HD3	2.17	0.45
1:A:212:LYS:HD3	6:A:965:HOH:O	2.16	0.45
1:A:148:GLN:HA	1:A:148:GLN:HE21	1.81	0.45
1:A:304:LYS:HG2	6:A:731:HOH:O	2.17	0.45
1:A:444:ILE:N	1:A:444:ILE:CD1	2.80	0.45
1:A:295:ASP:O	1:A:296:ASN:HB3	2.17	0.45
1:A:3:ASP:HA	1:A:548:GLN:HE21	1.82	0.45
1:A:82:ASN:HA	1:A:101:TRP:O	2.16	0.45
1:A:262:VAL:O	1:A:263:ASN:HB2	2.17	0.45
1:A:385:THR:O	1:A:389:GLN:HG3	2.17	0.45
1:A:408:SER:O	1:A:423:ARG:HA	2.17	0.45
1:A:54:TRP:HH2	1:A:108:THR:CG2	2.30	0.44
1:A:590:ARG:HB2	1:A:639:ASP:OD1	2.18	0.44
1:A:612:GLU:CD	1:A:643:PRO:HD3	2.42	0.44
1:A:82:ASN:OD1	1:A:99:GLY:HA2	2.17	0.44
1:A:194:LEU:HD13	2:B:3:GLC:H62	2.00	0.44
1:A:60:LYS:HA	1:A:60:LYS:HD2	1.65	0.44
1:A:178:ASN:O	1:A:192:LYS:HE2	2.18	0.44
1:A:563:ILE:HD13	1:A:579:PHE:HB2	1.99	0.44
1:A:417:ASP:O	1:A:436:ARG:HG3	2.18	0.44
1:A:630:VAL:HG12	1:A:631:TYR:CE1	2.52	0.44
1:A:651:LYS:HD2	1:A:667:ASN:OD1	2.17	0.43
1:A:627:ASN:HB2	1:A:633:TYR:N	2.33	0.43
1:A:40:ASP:HB2	1:A:48:LEU:HD23	1.99	0.43
1:A:159:ASP:HB2	1:A:164:LEU:HD21	1.99	0.43
1:A:212:LYS:HE2	1:A:242:PHE:HA	1.99	0.43
1:A:299:ASN:HB2	1:A:416:ASN:O	2.19	0.43
1:A:633:TYR:HA	1:A:635:ASN:N	2.33	0.43
1:A:16:VAL:HG21	1:A:395:ALA:HB1	2.00	0.43
1:A:74:ILE:HD12	1:A:76:ILE:HG23	2.00	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:666:SER:N	1:A:685:GLN:HE22	2.16	0.42
1:A:227:ARG:NH1	1:A:257:GLN:CB	2.79	0.42
1:A:496:THR:HG22	1:A:498:THR:O	2.18	0.42
1:A:83:ILE:HB	1:A:153:GLU:CD	2.45	0.42
1:A:637:TYR:O	1:A:638:TYR:HB2	2.18	0.42
1:A:650:PHE:O	1:A:668:HIS:HD2	2.02	0.42
1:A:466:GLY:HA3	1:A:488:ALA:HB2	2.00	0.42
1:A:290:ARG:NH2	1:A:325:ASP:CB	2.83	0.42
1:A:227:ARG:HG2	1:A:255:PHE:CE2	2.55	0.42
1:A:499:PRO:O	1:A:573:SER:HB2	2.20	0.42
1:A:563:ILE:CD1	1:A:579:PHE:CB	2.98	0.42
1:A:440:ALA:HA	1:A:441:PRO:HD3	1.73	0.42
1:A:538:GLY:O	1:A:541:ILE:HB	2.19	0.42
1:A:233:HIS:HB3	6:A:854:HOH:O	2.20	0.41
1:A:560:ASN:HA	1:A:579:PHE:O	2.19	0.41
1:A:589:VAL:HG13	1:A:678:ALA:HB3	2.02	0.41
1:A:33:ASN:HA	1:A:34:PRO:HD2	1.90	0.41
1:A:456:TYR:N	1:A:456:TYR:CD1	2.88	0.41
1:A:17:ILE:CG2	1:A:18:TYR:N	2.83	0.41
1:A:513:VAL:HB	1:A:553:ILE:HD12	2.02	0.41
1:A:627:ASN:CB	1:A:633:TYR:HB2	2.51	0.41
1:A:425:PHE:O	1:A:428:ASN:HB2	2.19	0.41
1:A:140:HIS:CD2	1:A:197:LEU:HD13	2.56	0.41
1:A:433:ALA:O	1:A:487:THR:HA	2.20	0.41
1:A:598:THR:HG22	1:A:602:GLN:HG3	2.03	0.41
2:B:2:GLC:C6	2:B:3:GLC:H3	2.46	0.41
1:A:45:ASN:HB3	1:A:48:LEU:HD22	2.01	0.41
1:A:402:PRO:HG2	1:A:428:ASN:ND2	2.36	0.41
1:A:648:ILE:O	1:A:669:THR:HA	2.21	0.41
1:A:140:HIS:NE2	1:A:197:LEU:HD13	2.36	0.41
1:A:633:TYR:CD1	1:A:633:TYR:C	2.99	0.41
1:A:42:THR:CG2	1:A:44:THR:CB	2.99	0.40
1:A:75:TRP:CD1	1:A:75:TRP:C	2.98	0.40
1:A:294:ARG:HB2	1:A:332:PHE:CE2	2.56	0.40
1:A:136:PHE:CD2	1:A:138:PRO:CG	3.00	0.40
1:A:512:GLY:O	1:A:552:LYS:HB3	2.22	0.40
1:A:562:ASN:HD22	1:A:575:VAL:CG1	2.24	0.40
1:A:631:TYR:HB2	1:A:637:TYR:HD1	1.86	0.40
1:A:451:LEU:HD22	1:A:456:TYR:CE2	2.56	0.40
1:A:544:TRP:CD1	1:A:544:TRP:C	2.99	0.40
1:A:365:TYR:CD1	1:A:365:TYR:N	2.89	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:562:ASN:HB3	1:A:575:VAL:HG13	2.03	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	684/686 (100%)	628 (92%)	52 (8%)	4 (1%)	21 18

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	483	ALA
1	A	629	VAL
1	A	46	LEU
1	A	148	GLN

5.3.2 Protein sidechains [i](#)

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
1	A	556/556 (100%)	486 (87%)	70 (13%)	4 2

All (70) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	42	THR
1	A	44	THR
1	A	48	LEU
1	A	67	THR
1	A	69	MET
1	A	71	VAL
1	A	75	TRP
1	A	80	VAL
1	A	82	ASN
1	A	90	SER
1	A	107	LYS
1	A	115	ILE
1	A	134	ILE
1	A	135	ASP
1	A	147	ASP
1	A	148	GLN
1	A	150	SER
1	A	163	LEU
1	A	169	ASN
1	A	173	ASN
1	A	181	THR
1	A	187	GLU
1	A	194	LEU
1	A	212	LYS
1	A	215	ILE
1	A	217	MET
1	A	227	ARG
1	A	232	LYS
1	A	279	SER
1	A	295	ASP
1	A	300	MET
1	A	304	LYS
1	A	331	ARG
1	A	335	SER
1	A	342	LEU
1	A	344	GLN
1	A	367	SER
1	A	375	ARG
1	A	384	SER
1	A	392	GLN
1	A	424	LYS
1	A	436	ARG
1	A	444	ILE

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Mol	Chain	Res	Type
1	A	449	THR
1	A	453	GLN
1	A	473	SER
1	A	491	GLN
1	A	520	ARG
1	A	524	SER
1	A	525	SER
1	A	534	THR
1	A	550	LYS
1	A	551	VAL
1	A	563	ILE
1	A	590	ARG
1	A	595	ASN
1	A	598	THR
1	A	600	LEU
1	A	602	GLN
1	A	606	LEU
1	A	609	SER
1	A	613	LEU
1	A	641	SER
1	A	646	LYS
1	A	658	SER
1	A	660	VAL
1	A	666	SER
1	A	669	THR
1	A	674	SER
1	A	675	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (19) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	19	GLN
1	A	62	ASN
1	A	88	ASN
1	A	119	GLN
1	A	120	ASN
1	A	139	ASN
1	A	269	ASN
1	A	336	ASN
1	A	410	GLN
1	A	416	ASN
1	A	453	GLN

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Mol	Chain	Res	Type
1	A	479	ASN
1	A	548	GLN
1	A	560	ASN
1	A	562	ASN
1	A	632	GLN
1	A	668	HIS
1	A	681	ASN
1	A	685	GLN

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 ⓘ

18 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	GLC	B	1	2	12,12,12	0.41	0	17,17,17	1.30	2 (11%)
2	GLC	B	2	2	11,11,12	0.47	0	15,15,17	1.28	2 (13%)
2	GLC	B	3	2	11,11,12	0.44	0	15,15,17	1.51	2 (13%)
2	GLC	B	4	2	11,11,12	0.48	0	15,15,17	1.85	1 (6%)
3	GLC	C	1	3	12,12,12	0.33	0	17,17,17	1.00	1 (5%)
3	GLC	C	2	3	11,11,12	0.53	0	15,15,17	1.52	3 (20%)
4	GLC	D	1	4	11,11,12	0.67	0	15,15,17	1.18	1 (6%)
4	GLC	D	2	4	11,11,12	0.50	0	15,15,17	1.50	1 (6%)
4	GLC	D	3	4	11,11,12	0.44	0	15,15,17	1.44	1 (6%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
4	GLC	D	4	4	11,11,12	0.28	0	15,15,17	1.04	1 (6%)
4	GLC	D	5	4	11,11,12	0.55	0	15,15,17	1.03	1 (6%)
4	GLC	D	6	4	11,11,12	0.70	0	15,15,17	1.04	1 (6%)
4	GLC	E	1	4	11,11,12	0.50	0	15,15,17	1.22	2 (13%)
4	GLC	E	2	4	11,11,12	0.32	0	15,15,17	0.80	0
4	GLC	E	3	4	11,11,12	0.52	0	15,15,17	0.94	0
4	GLC	E	4	4	11,11,12	0.63	0	15,15,17	1.26	1 (6%)
4	GLC	E	5	4	11,11,12	0.58	0	15,15,17	0.87	0
4	GLC	E	6	4	11,11,12	0.58	0	15,15,17	1.28	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	GLC	B	1	2	-	2/2/22/22	0/1/1/1
2	GLC	B	2	2	-	2/2/19/22	0/1/1/1
2	GLC	B	3	2	-	0/2/19/22	0/1/1/1
2	GLC	B	4	2	-	2/2/19/22	0/1/1/1
3	GLC	C	1	3	-	0/2/22/22	0/1/1/1
3	GLC	C	2	3	-	0/2/19/22	0/1/1/1
4	GLC	D	1	4	-	2/2/19/22	0/1/1/1
4	GLC	D	2	4	-	0/2/19/22	0/1/1/1
4	GLC	D	3	4	-	2/2/19/22	0/1/1/1
4	GLC	D	4	4	-	2/2/19/22	0/1/1/1
4	GLC	D	5	4	-	0/2/19/22	0/1/1/1
4	GLC	D	6	4	-	1/2/19/22	0/1/1/1
4	GLC	E	1	4	-	0/2/19/22	0/1/1/1
4	GLC	E	2	4	-	2/2/19/22	0/1/1/1
4	GLC	E	3	4	-	2/2/19/22	0/1/1/1
4	GLC	E	4	4	-	0/2/19/22	0/1/1/1
4	GLC	E	5	4	-	1/2/19/22	0/1/1/1
4	GLC	E	6	4	-	0/2/19/22	0/1/1/1

There are no bond length outliers.

All (22) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	4	GLC	C1-O5-C5	6.04	120.28	112.19
4	D	2	GLC	C1-O5-C5	4.73	118.53	112.19
2	B	3	GLC	C1-O5-C5	4.24	117.86	112.19
4	D	3	GLC	C1-O5-C5	4.13	117.72	112.19
2	B	1	GLC	C4-C3-C2	-3.30	105.03	110.83
3	C	2	GLC	O3-C3-C4	-3.05	103.19	110.38
2	B	2	GLC	C1-O5-C5	2.84	115.99	112.19
4	D	1	GLC	C1-O5-C5	2.83	115.97	112.19
2	B	3	GLC	C2-C3-C4	-2.68	106.15	110.86
4	E	6	GLC	C1-O5-C5	2.67	115.76	112.19
4	D	6	GLC	C1-O5-C5	2.61	115.69	112.19
4	E	1	GLC	O3-C3-C4	-2.43	104.65	110.38
4	E	6	GLC	C3-C4-C5	2.36	114.51	110.23
4	E	1	GLC	C3-C4-C5	2.35	114.49	110.23
4	D	5	GLC	O4-C4-C3	-2.18	105.25	110.38
2	B	2	GLC	O5-C1-C2	2.16	115.93	110.79
4	E	4	GLC	C1-C2-C3	-2.14	106.53	109.64
3	C	1	GLC	O4-C4-C3	-2.10	105.43	110.38
3	C	2	GLC	O3-C3-C2	-2.09	105.79	110.05
2	B	1	GLC	O5-C5-C4	-2.07	105.97	109.70
4	D	4	GLC	C1-O5-C5	2.05	114.94	112.19
3	C	2	GLC	C6-C5-C4	-2.05	107.99	113.02

There are no chirality outliers.

All (18) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	E	2	GLC	O5-C5-C6-O6
2	B	2	GLC	O5-C5-C6-O6
4	D	3	GLC	C4-C5-C6-O6
2	B	2	GLC	C4-C5-C6-O6
4	E	2	GLC	C4-C5-C6-O6
2	B	1	GLC	O5-C5-C6-O6
2	B	4	GLC	C4-C5-C6-O6
4	D	3	GLC	O5-C5-C6-O6
2	B	1	GLC	C4-C5-C6-O6
2	B	4	GLC	O5-C5-C6-O6
4	D	1	GLC	C4-C5-C6-O6
4	D	4	GLC	O5-C5-C6-O6
4	D	4	GLC	C4-C5-C6-O6
4	D	1	GLC	O5-C5-C6-O6
4	E	3	GLC	O5-C5-C6-O6
4	E	3	GLC	C4-C5-C6-O6

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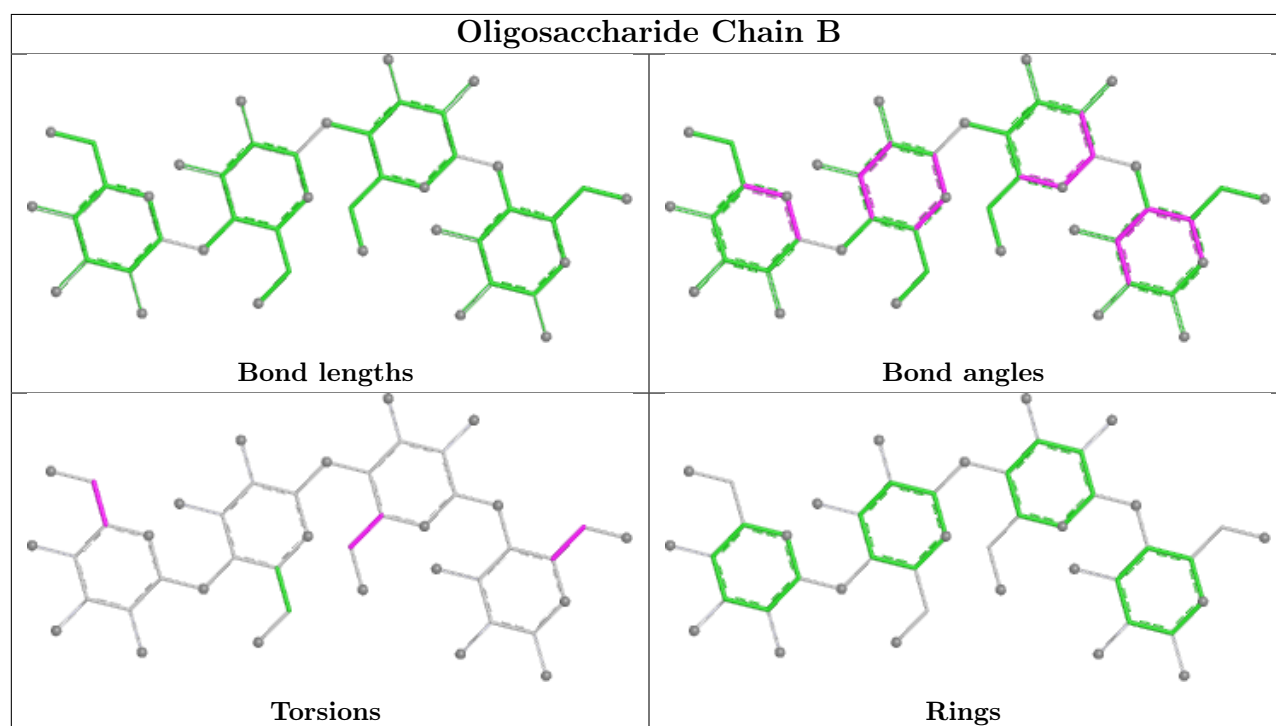
Mol	Chain	Res	Type	Atoms
4	E	5	GLC	O5-C5-C6-O6
4	D	6	GLC	O5-C5-C6-O6

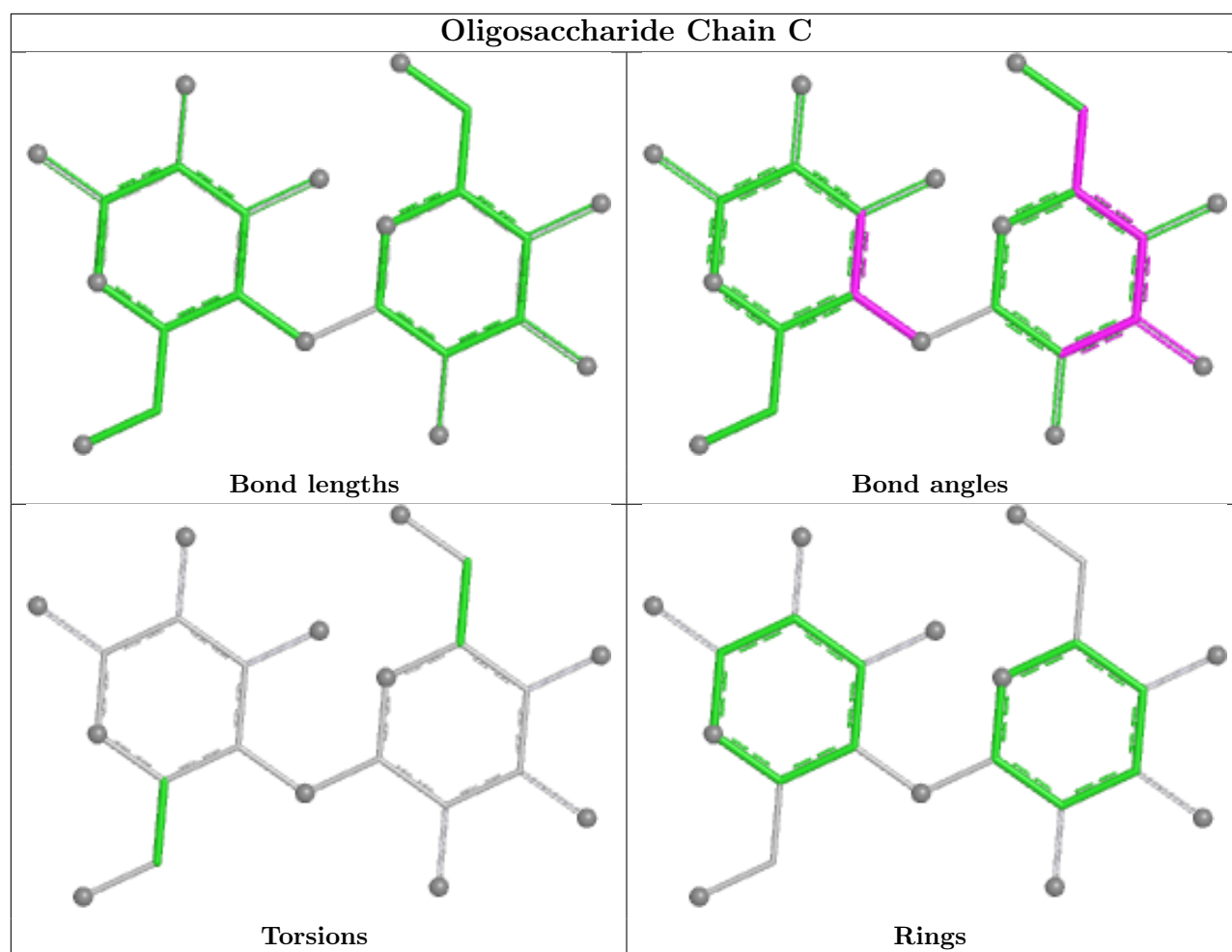
There are no ring outliers.

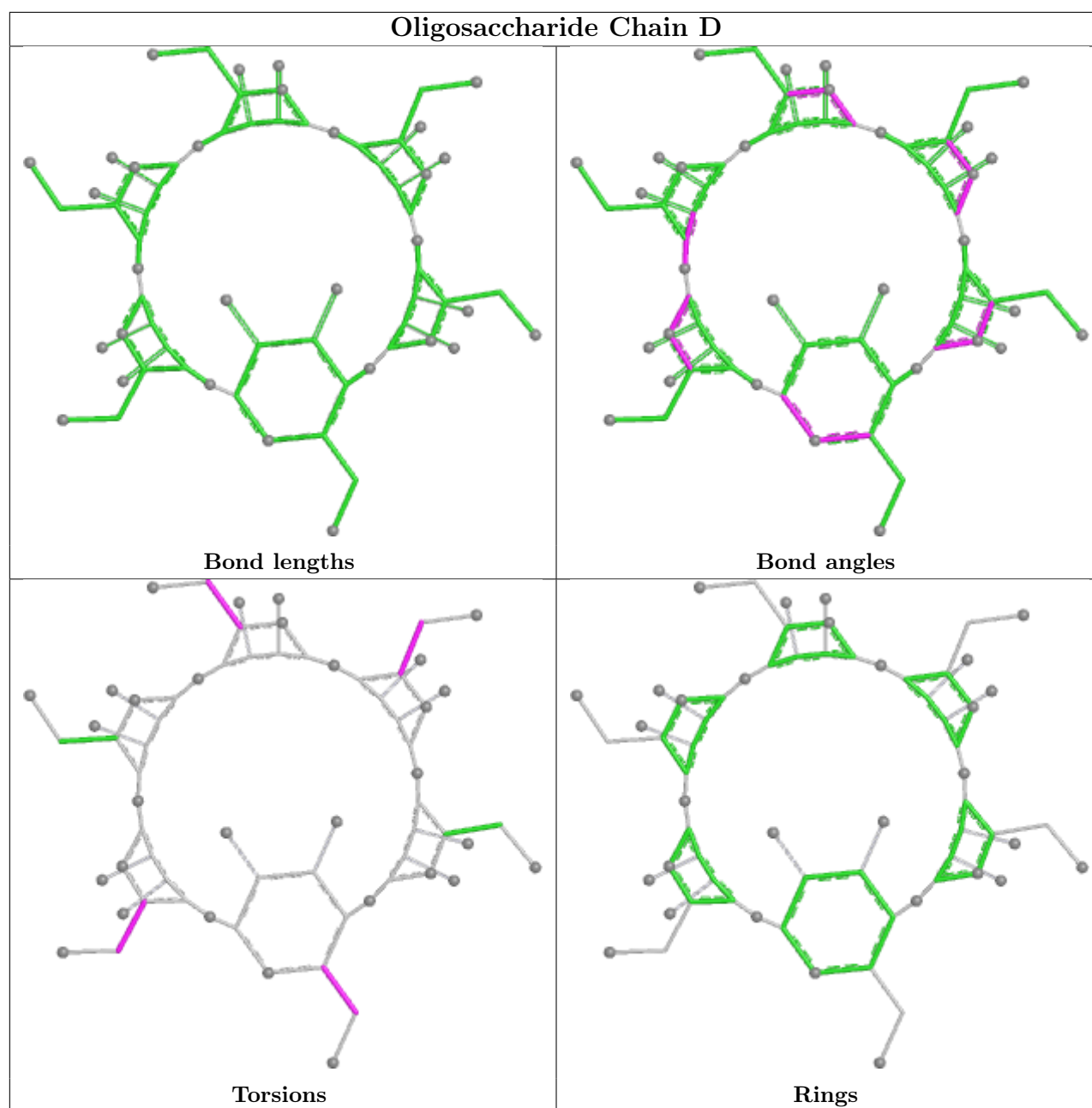
4 monomers are involved in 6 short contacts:

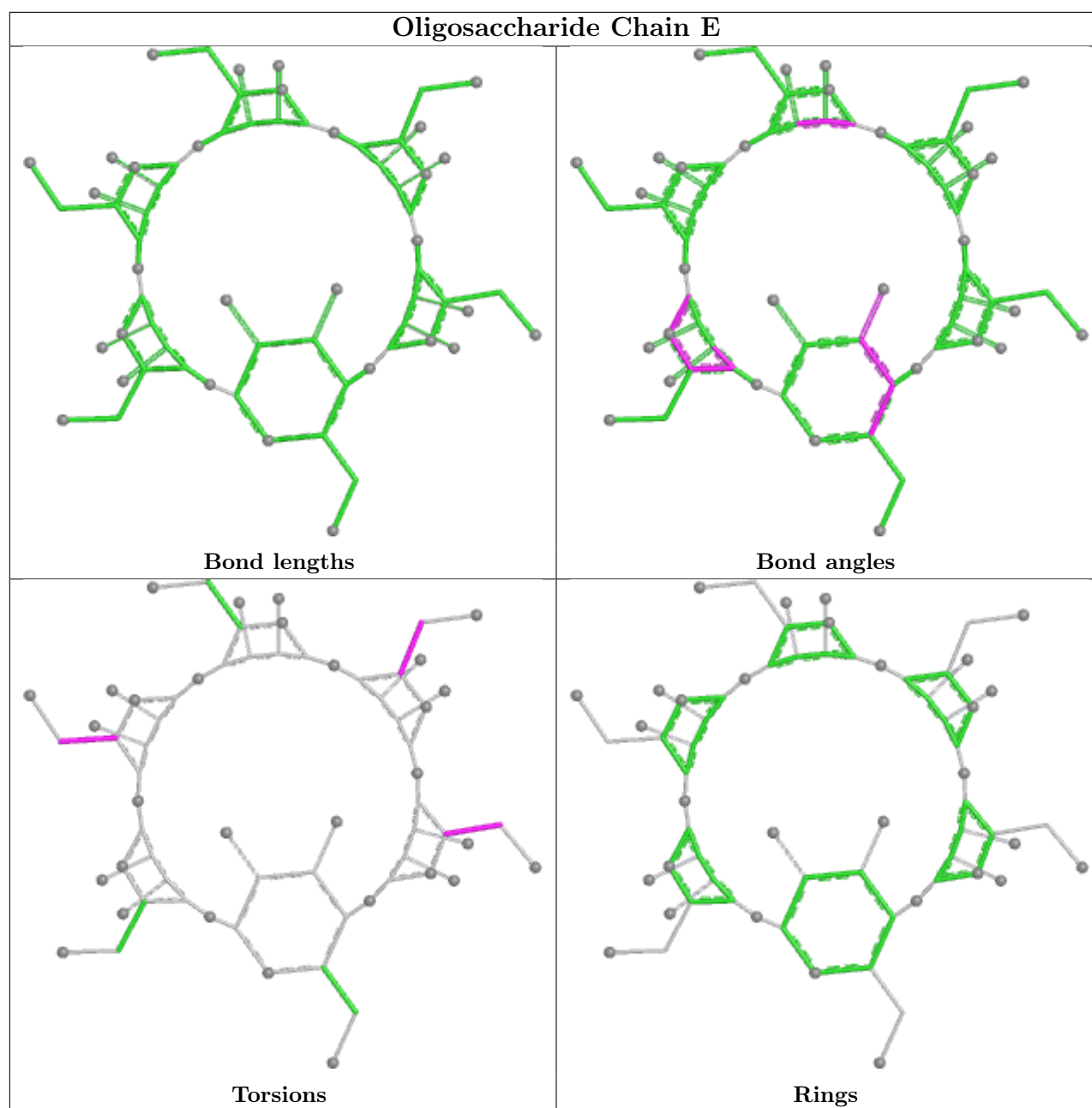
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	4	GLC	1	0
2	B	3	GLC	4	0
4	E	4	GLC	1	0
2	B	2	GLC	3	0

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 2 ligands modelled in this entry, 2 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

There are no such residues in this entry.

5.8 Polymer linkage issues

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	686/686 (100%)	-0.11	13 (1%) 66 69	2, 9, 22, 42	0

All (13) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	73	ALA	3.0
1	A	466	GLY	2.9
1	A	89	TYR	2.4
1	A	599	ALA	2.4
1	A	197	LEU	2.3
1	A	145	SER	2.3
1	A	127	ALA	2.3
1	A	494	ALA	2.2
1	A	41	GLY	2.2
1	A	387	ALA	2.2
1	A	121	LEU	2.1
1	A	254	THR	2.0
1	A	598	THR	2.0

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

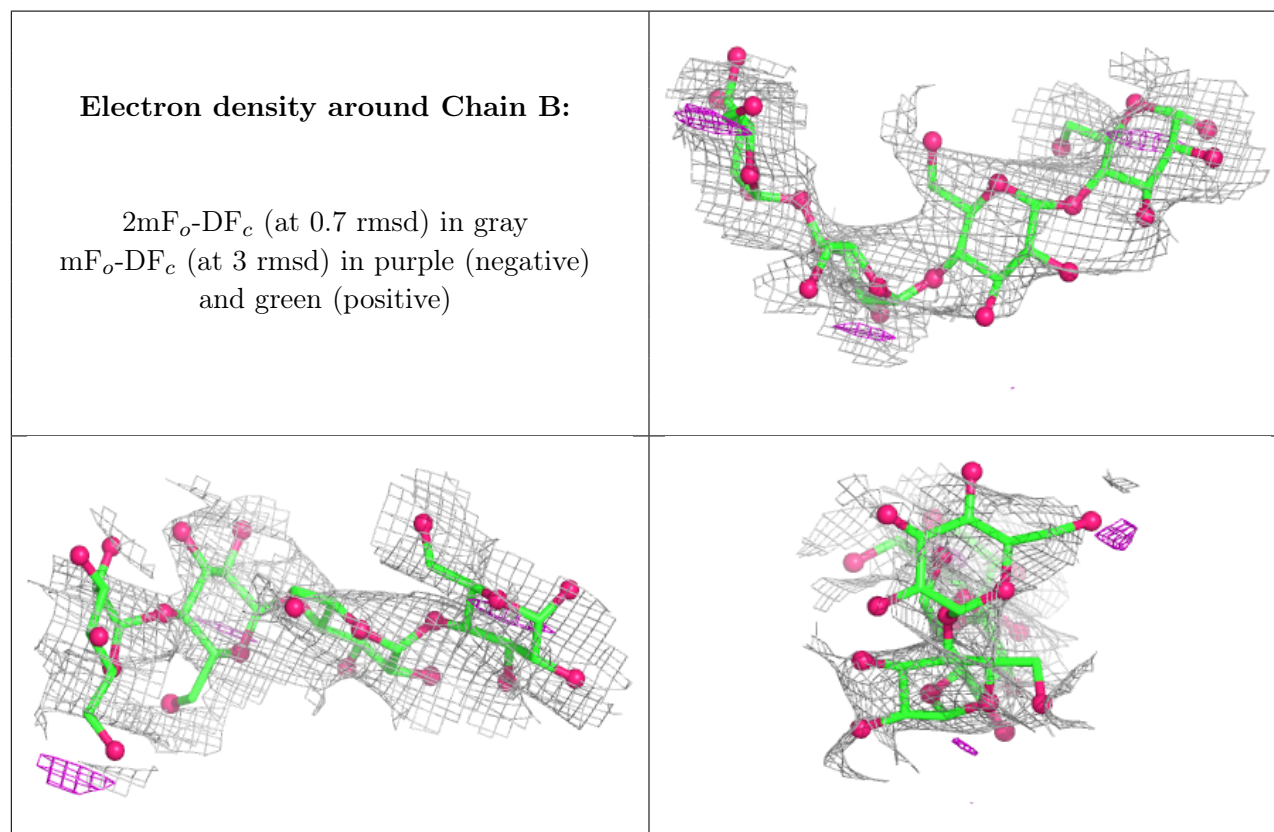
There are no non-standard protein/DNA/RNA residues in this entry.

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.

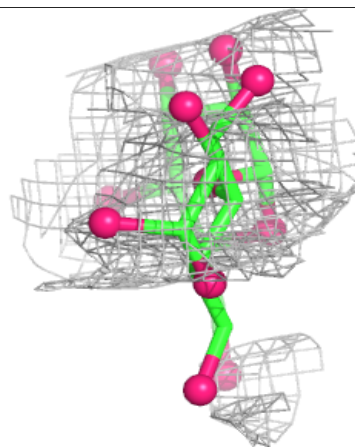
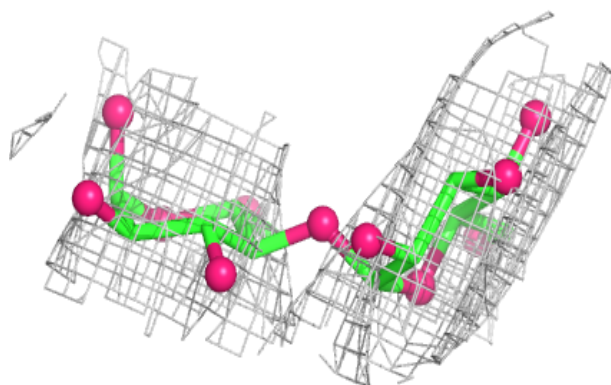
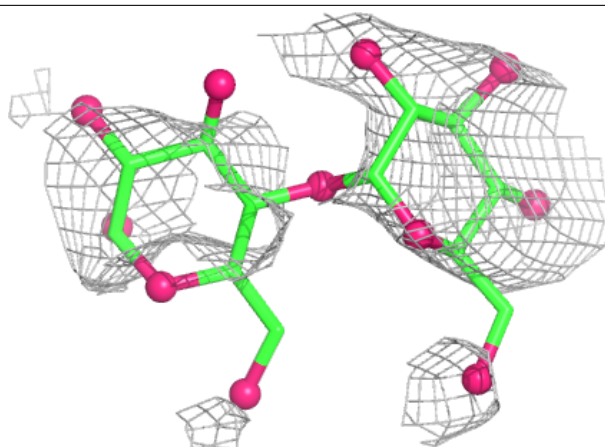
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	GLC	B	1	12/12	0.80	0.10	56,56,61,66	0
4	GLC	E	1	11/12	0.83	0.12	45,45,51,56	0
4	GLC	E	4	11/12	0.83	0.14	44,44,49,54	0
4	GLC	D	2	11/12	0.84	0.11	40,41,46,51	0
4	GLC	D	3	11/12	0.88	0.10	40,40,45,50	0
2	GLC	B	3	11/12	0.89	0.12	54,55,59,64	0
3	GLC	C	1	12/12	0.89	0.12	3,3,8,14	0
4	GLC	E	3	11/12	0.90	0.07	45,45,50,55	0
4	GLC	D	5	11/12	0.90	0.09	33,33,38,42	0
4	GLC	D	6	11/12	0.91	0.12	32,33,38,42	0
2	GLC	B	4	11/12	0.92	0.11	52,52,57,62	0
4	GLC	E	2	11/12	0.92	0.07	46,46,51,56	0
4	GLC	D	4	11/12	0.93	0.07	37,38,42,47	0
2	GLC	B	2	11/12	0.94	0.09	55,55,59,64	0
3	GLC	C	2	11/12	0.94	0.06	4,4,9,14	0
4	GLC	D	1	11/12	0.94	0.07	36,37,42,47	0
4	GLC	E	6	11/12	0.95	0.07	43,43,48,53	0
4	GLC	E	5	11/12	0.97	0.07	43,44,48,53	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.



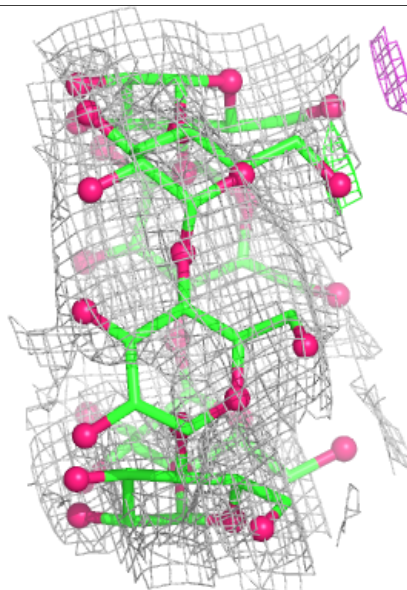
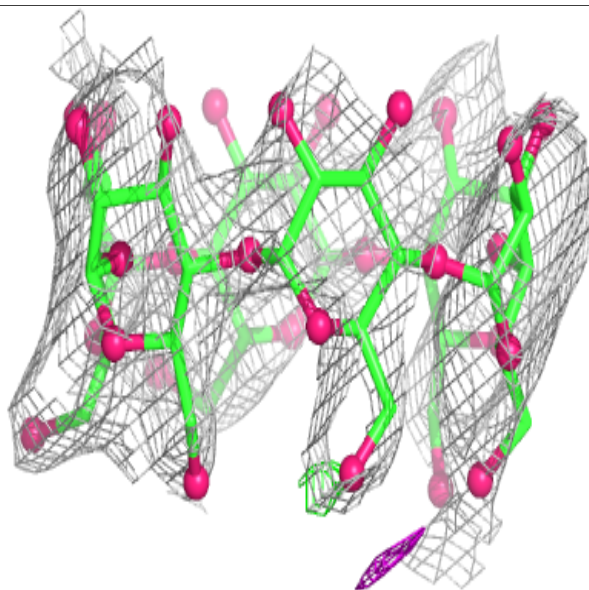
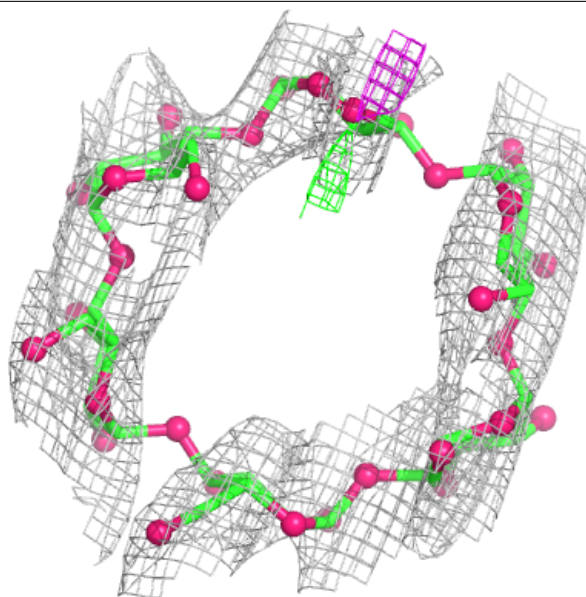
Electron density around Chain C:

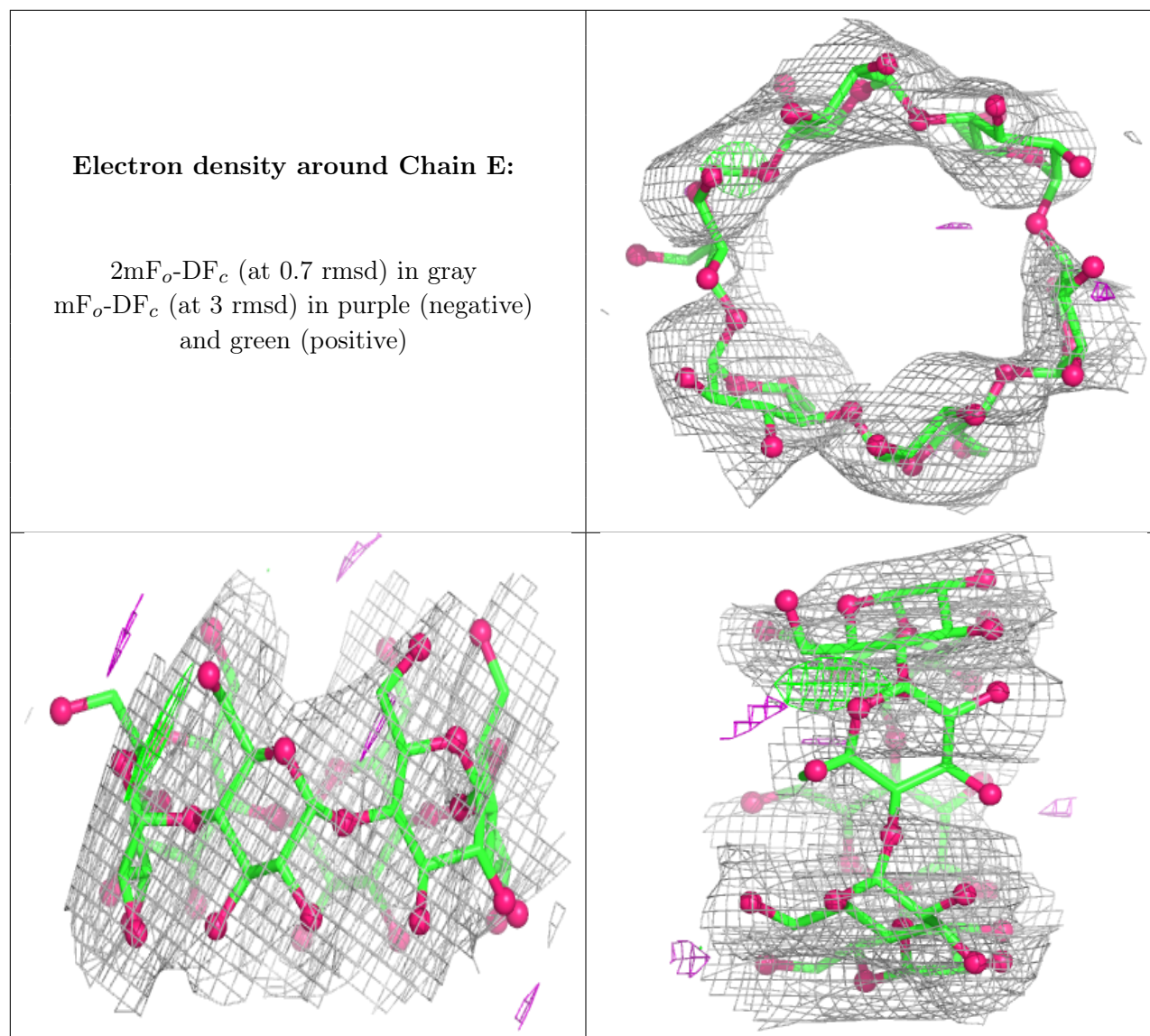
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around Chain D:

$2mF_o-DF_c$ (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(Å ²)	Q<0.9
5	CA	A	690	1/1	0.93	0.04	12,12,12,12	0
5	CA	A	691	1/1	0.94	0.10	17,17,17,17	0

6.5 Other polymers [i](#)

There are no such residues in this entry.