



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 02:21 AM UTC

PDB ID : 1CGW / pdb_00001cgw
Title : SITE DIRECTED MUTATIONS OF THE ACTIVE SITE RESIDUE TYROSINE 195 OF CYCLODEXTRIN GLYCOSYLTRANSFERASE FROM BACILLUS CIRCULANS STRAIN 251 AFFECTING ACTIVITY AND PRODUCT SPECIFICITY
Authors : Strokopytov, B.V.; Dijkstra, B.W.
Deposited on : 1994-08-05
Resolution : 2.50 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

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

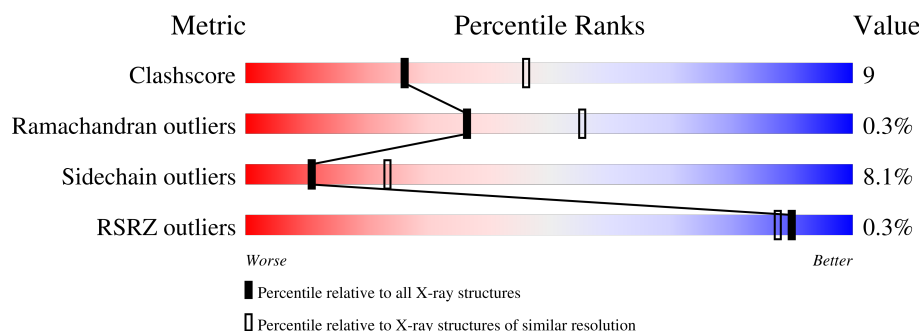
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

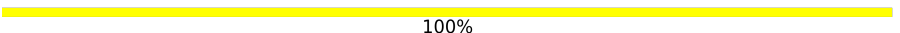
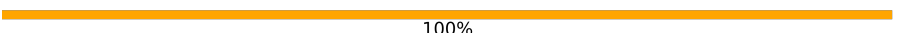
The reported resolution of this entry is 2.50 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



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

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	686	 72% 24% . .
2	B	2	 50% 50%
2	C	2	 100%
2	D	2	 100%

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 5506 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 CYCLOMALTODEXTRIN GLUCANOTRANSFERASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	686	Total	C	N	O	S	0	0	0
			5256	3314	900	1026	16			

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	195	GLY	TYR	engineered mutation	UNP P43379

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



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

- Molecule 3 is CALCIUM ION (CCD ID: CA) (formula: Ca).

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

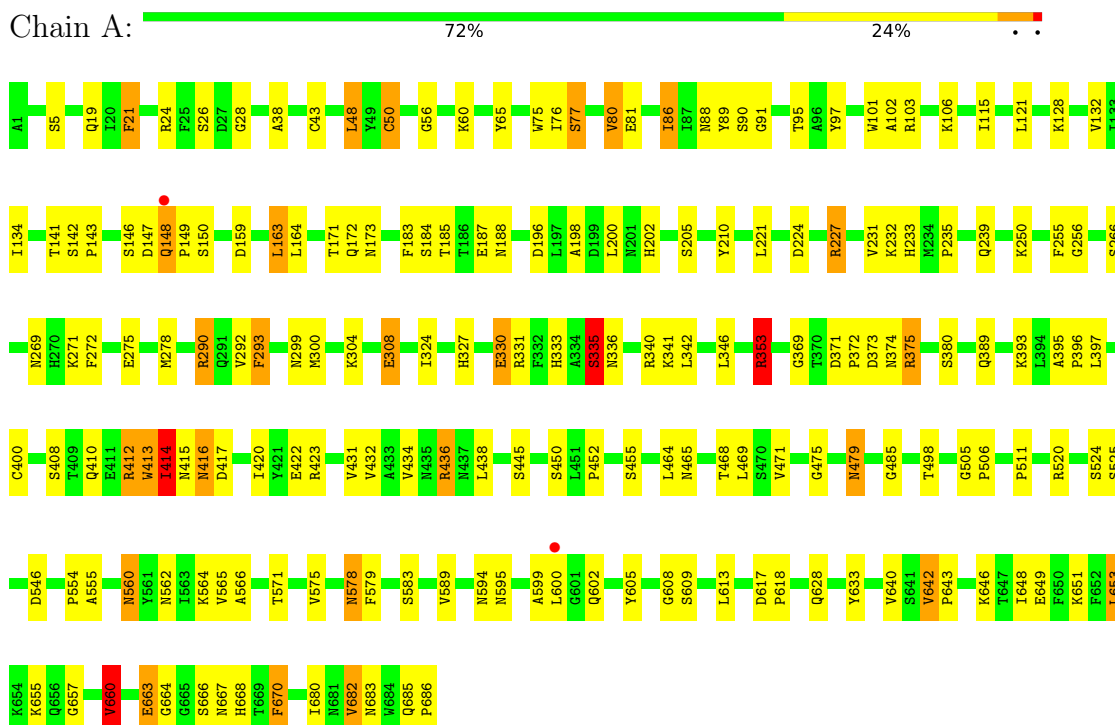
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	179	Total 179	O 179	0	0

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: CYCLOMALTODEXTRIN GLUCANOTRANSFERASE



• Molecule 2: alpha-D-glucopyranose-(1-4)-alpha-D-glucopyranose



• Molecule 2: alpha-D-glucopyranose-(1-4)-alpha-D-glucopyranose



• Molecule 2: alpha-D-glucopyranose-(1-4)-alpha-D-glucopyranose

Chain D:

100%

GLU1
GLU2

4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	121.09Å 111.18Å 66.80Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.00 – 2.50 20.00 – 2.55	Depositor EDS
% Data completeness (in resolution range)	(Not available) (20.00-2.50) 91.8 (20.00-2.55)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.73 (at 2.54Å)	Xtriage
Refinement program	PROLSQ, TNT	Depositor
R, R_{free}	0.160 , (Not available) 0.146 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å ²)	24.8	Xtriage
Anisotropy	0.070	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 83.9	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	0.96	EDS
Total number of atoms	5506	wwPDB-VP
Average B, all atoms (Å ²)	24.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.72% 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	1.01	2/5385 (0.0%)	1.59	65/7339 (0.9%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	413	TRP	C-N	-5.56	1.30	1.33
1	A	290	ARG	CZ-NH2	-5.00	1.26	1.33

All (65) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	373	ASP	N-CA-C	9.57	124.00	111.75
1	A	414	ILE	CA-C-O	7.92	125.96	120.66
1	A	336	ASN	N-CA-CB	-7.89	98.18	111.20
1	A	353	ARG	CD-NE-CZ	7.74	135.23	124.40
1	A	609	SER	N-CA-C	7.74	122.15	112.87
1	A	660	VAL	N-CA-C	7.42	120.17	108.89
1	A	266	SER	CA-C-N	7.32	126.95	119.19
1	A	266	SER	C-N-CA	7.32	126.95	119.19
1	A	412	ARG	N-CA-CB	7.29	120.94	110.88
1	A	353	ARG	NE-CZ-NH2	-7.22	112.70	119.20
1	A	224	ASP	CA-CB-CG	-7.21	105.39	112.60
1	A	413	TRP	CA-C-N	-7.11	114.86	121.97
1	A	413	TRP	C-N-CA	-7.11	114.86	121.97
1	A	80	VAL	N-CA-C	6.96	118.08	110.21
1	A	663	GLU	N-CA-C	-6.96	100.68	110.50
1	A	668	HIS	N-CA-C	-6.88	100.61	110.59
1	A	331	ARG	NE-CZ-NH2	-6.80	113.08	119.20
1	A	414	ILE	N-CA-C	6.73	115.91	107.70
1	A	50	CYS	CB-CA-C	6.48	119.95	109.07

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	335	SER	CA-C-N	6.35	132.46	122.49
1	A	335	SER	C-N-CA	6.35	132.46	122.49
1	A	642	VAL	O-C-N	6.33	126.15	121.72
1	A	327	HIS	CA-CB-CG	-6.33	107.47	113.80
1	A	233	HIS	CA-CB-CG	-6.30	107.50	113.80
1	A	416	ASN	CB-CA-C	-6.27	100.19	110.85
1	A	184	SER	N-CA-C	6.16	118.07	111.36
1	A	292	VAL	N-CA-C	6.10	116.26	110.53
1	A	28	GLY	CA-C-N	-6.02	116.36	122.74
1	A	28	GLY	C-N-CA	-6.02	116.36	122.74
1	A	682	VAL	N-CA-C	6.00	115.84	108.53
1	A	266	SER	O-C-N	5.91	126.65	121.39
1	A	21	PHE	CA-CB-CG	-5.86	107.94	113.80
1	A	608	GLY	CA-C-N	5.86	131.88	120.99
1	A	608	GLY	C-N-CA	5.86	131.88	120.99
1	A	657	GLY	N-CA-C	-5.85	100.09	110.86
1	A	600	LEU	N-CA-C	5.71	118.61	110.10
1	A	205	SER	N-CA-C	5.70	117.49	111.28
1	A	546	ASP	CA-CB-CG	-5.61	107.00	112.60
1	A	485	GLY	CA-C-N	5.52	125.94	120.31
1	A	485	GLY	C-N-CA	5.52	125.94	120.31
1	A	335	SER	N-CA-C	5.42	122.35	110.80
1	A	102	ALA	N-CA-C	5.38	118.04	110.23
1	A	202	HIS	CA-CB-CG	-5.38	108.42	113.80
1	A	410	GLN	N-CA-C	5.37	117.17	108.26
1	A	290	ARG	NE-CZ-NH1	5.35	126.85	121.50
1	A	479	ASN	CA-CB-CG	-5.33	107.27	112.60
1	A	554	PRO	CB-CA-C	-5.32	104.26	111.12
1	A	196	ASP	CB-CG-OD2	-5.29	106.23	118.40
1	A	65	TYR	N-CA-C	5.27	117.45	111.02
1	A	452	PRO	N-CA-CB	5.18	108.15	103.33
1	A	346	LEU	CB-CA-C	-5.18	102.72	110.90
1	A	640	VAL	N-CA-C	5.18	116.17	108.46
1	A	648	ILE	N-CA-C	5.16	116.15	108.46
1	A	86	ILE	N-CA-C	-5.14	102.10	108.89
1	A	76	ILE	N-CA-C	5.12	116.98	108.99
1	A	389	GLN	CB-CG-CD	-5.12	103.90	112.60
1	A	293	PHE	CA-CB-CG	5.09	118.89	113.80
1	A	172	GLN	N-CA-C	-5.07	105.88	111.71
1	A	148	GLN	CA-C-N	5.07	126.17	119.84
1	A	148	GLN	C-N-CA	5.07	126.17	119.84
1	A	221	LEU	N-CA-CB	5.04	118.89	110.41

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	132	VAL	N-CA-C	5.03	115.15	108.11
1	A	578	ASN	CA-CB-CG	5.03	117.63	112.60
1	A	101	TRP	CA-C-N	5.02	127.83	120.71
1	A	101	TRP	C-N-CA	5.02	127.83	120.71

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	5256	0	5019	97	0
2	B	23	0	21	0	0
2	C	23	0	21	0	0
2	D	23	0	21	1	0
3	A	2	0	0	0	0
4	A	179	0	0	3	0
All	All	5506	0	5082	98	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 9.

All (98) 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:88:ASN:HD21	1:A:91:GLY:H	1.01	0.92
1:A:142:SER:HB2	1:A:143:PRO:HD2	1.63	0.81
1:A:88:ASN:HD21	1:A:91:GLY:N	1.80	0.78
1:A:159:ASP:HB2	1:A:164:LEU:HD11	1.68	0.74
1:A:560:ASN:HD21	1:A:578:ASN:HA	1.53	0.74
1:A:643:PRO:HB2	1:A:646:LYS:HG3	1.68	0.74
1:A:594:ASN:HB2	1:A:683:ASN:OD1	1.91	0.70
1:A:147:ASP:C	1:A:149:PRO:HD3	2.16	0.70
1:A:163:LEU:HD22	1:A:164:LEU:N	2.07	0.69
1:A:88:ASN:ND2	1:A:91:GLY:H	1.86	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:333:HIS:HB2	1:A:342:LEU:HD13	1.79	0.64
1:A:304:LYS:O	1:A:308:GLU:HG2	1.98	0.64
1:A:670:PHE:CE2	1:A:680:ILE:HD11	2.33	0.63
1:A:562:ASN:HB3	1:A:575:VAL:HG13	1.82	0.61
1:A:395:ALA:HB3	1:A:396:PRO:HD3	1.82	0.61
1:A:227:ARG:HG2	1:A:255:PHE:CE2	2.36	0.60
1:A:589:VAL:HG11	1:A:680:ILE:HD11	1.83	0.59
1:A:560:ASN:ND2	1:A:578:ASN:HA	2.15	0.59
1:A:81:GLU:OE1	1:A:103:ARG:HD3	2.02	0.59
1:A:471:VAL:CG1	1:A:475:GLY:HA2	2.32	0.59
1:A:653:LEU:HD12	1:A:660:VAL:CG2	2.34	0.58
1:A:142:SER:HB2	1:A:143:PRO:CD	2.33	0.57
1:A:445:SER:HB3	1:A:479:ASN:ND2	2.21	0.55
1:A:141:THR:OG1	1:A:198:ALA:HB3	2.07	0.55
1:A:670:PHE:CZ	1:A:680:ILE:HD11	2.41	0.54
1:A:653:LEU:HD23	1:A:653:LEU:N	2.22	0.54
1:A:121:LEU:HD12	1:A:121:LEU:O	2.07	0.54
1:A:422:GLU:HG3	1:A:431:VAL:HG22	1.90	0.54
1:A:353:ARG:HD2	4:A:1044:HOH:O	2.08	0.53
1:A:38:ALA:O	1:A:48:LEU:HB3	2.09	0.52
1:A:185:THR:OG1	1:A:188:ASN:HB2	2.09	0.52
1:A:520:ARG:HG2	1:A:520:ARG:HH11	1.75	0.51
1:A:187:GLU:CD	1:A:628:GLN:HG2	2.36	0.51
1:A:148:GLN:N	1:A:149:PRO:HD3	2.23	0.51
1:A:232:LYS:O	1:A:232:LYS:HG2	2.10	0.51
1:A:642:VAL:HB	1:A:643:PRO:HD2	1.93	0.51
1:A:589:VAL:HG11	1:A:680:ILE:CD1	2.41	0.50
1:A:599:ALA:O	1:A:602:GLN:HB3	2.12	0.50
1:A:663:GLU:HG2	1:A:685:GLN:O	2.12	0.50
1:A:566:ALA:HA	1:A:571:THR:O	2.11	0.50
1:A:417:ASP:O	1:A:436:ARG:HG3	2.11	0.50
1:A:505:GLY:HA2	1:A:506:PRO:C	2.38	0.49
1:A:414:ILE:HG12	1:A:415:ASN:N	2.26	0.49
1:A:232:LYS:HB3	4:A:1166:HOH:O	2.12	0.49
1:A:397:LEU:HA	1:A:400:CYS:SG	2.53	0.48
1:A:649:GLU:OE1	1:A:667:ASN:HB2	2.13	0.48
1:A:271:LYS:O	1:A:275:GLU:HG3	2.14	0.48
1:A:333:HIS:HB2	1:A:342:LEU:CD1	2.44	0.47
1:A:408:SER:O	1:A:423:ARG:HA	2.14	0.47
1:A:468:THR:HG22	1:A:469:LEU:N	2.29	0.47
1:A:562:ASN:HB3	1:A:575:VAL:CG1	2.44	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:445:SER:CB	1:A:479:ASN:ND2	2.78	0.47
1:A:300:MET:HB2	1:A:415:ASN:O	2.14	0.47
1:A:420:ILE:HA	1:A:432:VAL:O	2.14	0.47
1:A:19:GLN:HG3	1:A:75:TRP:CD2	2.49	0.47
1:A:605:TYR:CE1	1:A:655:LYS:HB2	2.50	0.47
1:A:664:GLY:O	1:A:685:GLN:HB2	2.15	0.47
1:A:163:LEU:HD22	1:A:163:LEU:C	2.38	0.46
1:A:511:PRO:HB3	1:A:555:ALA:HA	1.98	0.46
1:A:653:LEU:HD12	1:A:660:VAL:HG21	1.97	0.46
1:A:293:PHE:HE1	1:A:434:VAL:HG11	1.80	0.46
1:A:235:PRO:O	1:A:239:GLN:HG3	2.16	0.46
1:A:340:ARG:O	1:A:340:ARG:HG3	2.15	0.46
1:A:617:ASP:OD1	1:A:618:PRO:HD2	2.16	0.46
1:A:272:PHE:CZ	1:A:278:MET:HE3	2.51	0.45
1:A:506:PRO:O	1:A:579:PHE:HE1	2.00	0.45
1:A:227:ARG:CG	1:A:255:PHE:CE2	3.00	0.45
1:A:148:GLN:N	1:A:149:PRO:CD	2.79	0.45
1:A:374:ASN:OD1	1:A:375:ARG:HD3	2.17	0.45
1:A:24:ARG:HD2	1:A:375:ARG:O	2.17	0.44
1:A:77:SER:HB3	4:A:1021:HOH:O	2.17	0.44
1:A:48:LEU:HD12	1:A:95:THR:HG23	1.99	0.44
1:A:121:LEU:HD12	1:A:121:LEU:C	2.43	0.44
1:A:330:GLU:HB3	1:A:369:GLY:HA2	2.00	0.44
1:A:272:PHE:HZ	1:A:278:MET:HE3	1.82	0.44
1:A:106:LYS:HD2	1:A:210:TYR:HE1	1.83	0.43
2:D:1:GLC:H62	2:D:2:GLC:C5	2.48	0.43
1:A:371:ASP:OD1	1:A:372:PRO:HA	2.17	0.43
1:A:685:GLN:HA	1:A:686:PRO:HD3	1.89	0.43
1:A:227:ARG:HD2	1:A:227:ARG:C	2.44	0.42
1:A:464:LEU:O	1:A:465:ASN:HB2	2.20	0.42
1:A:290:ARG:NH2	1:A:330:GLU:O	2.53	0.42
1:A:395:ALA:N	1:A:396:PRO:CD	2.82	0.42
1:A:106:LYS:HD2	1:A:210:TYR:CE1	2.55	0.42
1:A:227:ARG:HG2	1:A:255:PHE:CD2	2.54	0.42
1:A:413:TRP:O	1:A:414:ILE:HB	2.19	0.42
1:A:89:TYR:O	1:A:90:SER:HB2	2.18	0.41
1:A:43:CYS:HB2	1:A:50:CYS:HB3	1.86	0.41
1:A:231:VAL:HG11	1:A:256:GLY:HA3	2.01	0.41
1:A:60:LYS:HD2	1:A:60:LYS:HA	1.70	0.41
1:A:299:ASN:HB2	1:A:416:ASN:O	2.21	0.41
1:A:134:ILE:O	1:A:134:ILE:HG13	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:417:ASP:OD1	1:A:436:ARG:HD2	2.21	0.41
1:A:583:SER:HB2	1:A:643:PRO:HG3	2.03	0.41
1:A:633:TYR:CD1	1:A:633:TYR:C	2.98	0.41
1:A:97:TYR:CD2	1:A:97:TYR:C	2.99	0.40
1:A:26:SER:O	1:A:56:GLY:HA3	2.21	0.40
1:A:200:LEU:HD23	1:A:200:LEU:HA	1.92	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%)	645 (94%)	37 (5%)	2 (0%)	36 55

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	335	SER
1	A	183	PHE

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	555/555 (100%)	510 (92%)	45 (8%)	11 23

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

Mol	Chain	Res	Type
1	A	5	SER
1	A	21	PHE
1	A	48	LEU
1	A	77	SER
1	A	80	VAL
1	A	86	ILE
1	A	115	ILE
1	A	128	LYS
1	A	146	SER
1	A	150	SER
1	A	163	LEU
1	A	171	THR
1	A	173	ASN
1	A	227	ARG
1	A	250	LYS
1	A	269	ASN
1	A	308	GLU
1	A	324	ILE
1	A	330	GLU
1	A	335	SER
1	A	341	LYS
1	A	353	ARG
1	A	375	ARG
1	A	380	SER
1	A	393	LYS
1	A	412	ARG
1	A	414	ILE
1	A	436	ARG
1	A	438	LEU
1	A	450	SER
1	A	455	SER
1	A	498	THR
1	A	524	SER
1	A	525	SER
1	A	560	ASN
1	A	564	LYS
1	A	565	VAL
1	A	595	ASN
1	A	613	LEU
1	A	651	LYS
1	A	653	LEU
1	A	660	VAL

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Mol	Chain	Res	Type
1	A	666	SER
1	A	670	PHE
1	A	682	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (15) 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	169	ASN
1	A	269	ASN
1	A	389	GLN
1	A	392	GLN
1	A	410	GLN
1	A	416	ASN
1	A	479	ASN
1	A	560	ASN
1	A	632	GLN
1	A	685	GLN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

6 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	0.38	0	17,17,17	0.98	0
2	GLC	B	2	2	11,11,12	0.39	0	15,15,17	1.64	2 (13%)
2	GLC	C	1	2	12,12,12	0.36	0	17,17,17	1.51	3 (17%)
2	GLC	C	2	2	11,11,12	0.46	0	15,15,17	1.30	3 (20%)
2	GLC	D	1	2	12,12,12	0.31	0	17,17,17	1.28	2 (11%)
2	GLC	D	2	2	11,11,12	0.30	0	15,15,17	1.56	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	-	0/2/22/22	0/1/1/1
2	GLC	B	2	2	-	2/2/19/22	0/1/1/1
2	GLC	C	1	2	-	2/2/22/22	0/1/1/1
2	GLC	C	2	2	-	0/2/19/22	0/1/1/1
2	GLC	D	1	2	-	0/2/22/22	0/1/1/1
2	GLC	D	2	2	-	2/2/19/22	0/1/1/1

There are no bond length outliers.

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	2	GLC	C1-O5-C5	4.95	118.82	112.19
2	B	2	GLC	C1-C2-C3	-3.94	103.91	109.64
2	C	1	GLC	O5-C1-C2	3.42	116.31	110.30
2	C	2	GLC	O2-C2-C3	-2.76	104.43	110.15
2	D	1	GLC	O4-C4-C3	-2.64	104.15	110.38
2	C	2	GLC	O5-C5-C6	-2.37	103.05	107.66
2	D	1	GLC	O5-C1-C2	2.37	114.46	110.30
2	B	2	GLC	O2-C2-C3	-2.33	105.33	110.15
2	D	2	GLC	O5-C1-C2	2.30	116.27	110.79
2	C	1	GLC	C4-C3-C2	-2.27	106.85	110.83
2	C	2	GLC	C1-O5-C5	2.24	115.19	112.19
2	C	1	GLC	O4-C4-C5	-2.08	104.19	109.32

There are no chirality outliers.

All (6) torsion outliers are listed below:

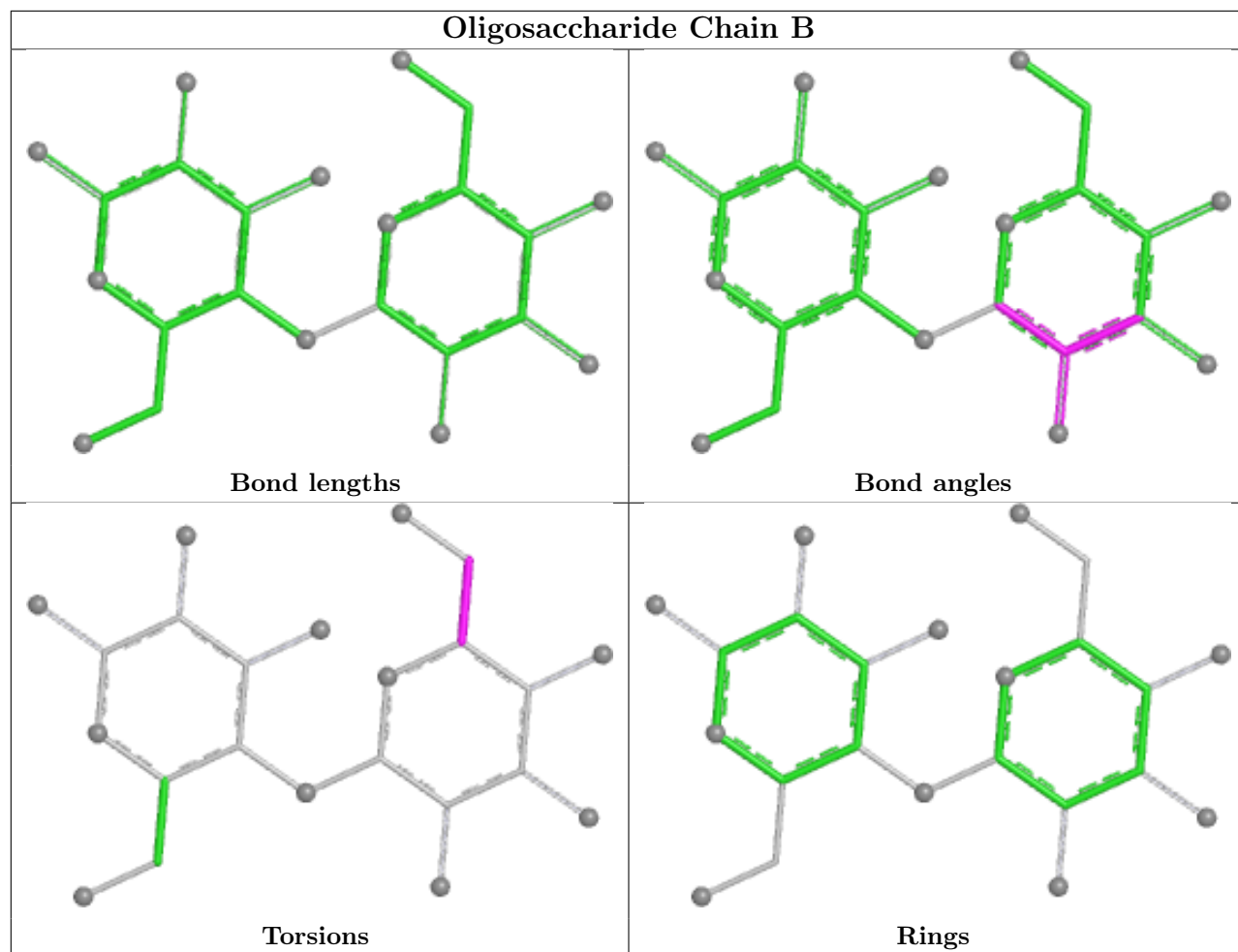
Mol	Chain	Res	Type	Atoms
2	D	2	GLC	C4-C5-C6-O6
2	D	2	GLC	O5-C5-C6-O6
2	C	1	GLC	O5-C5-C6-O6
2	C	1	GLC	C4-C5-C6-O6
2	B	2	GLC	C4-C5-C6-O6
2	B	2	GLC	O5-C5-C6-O6

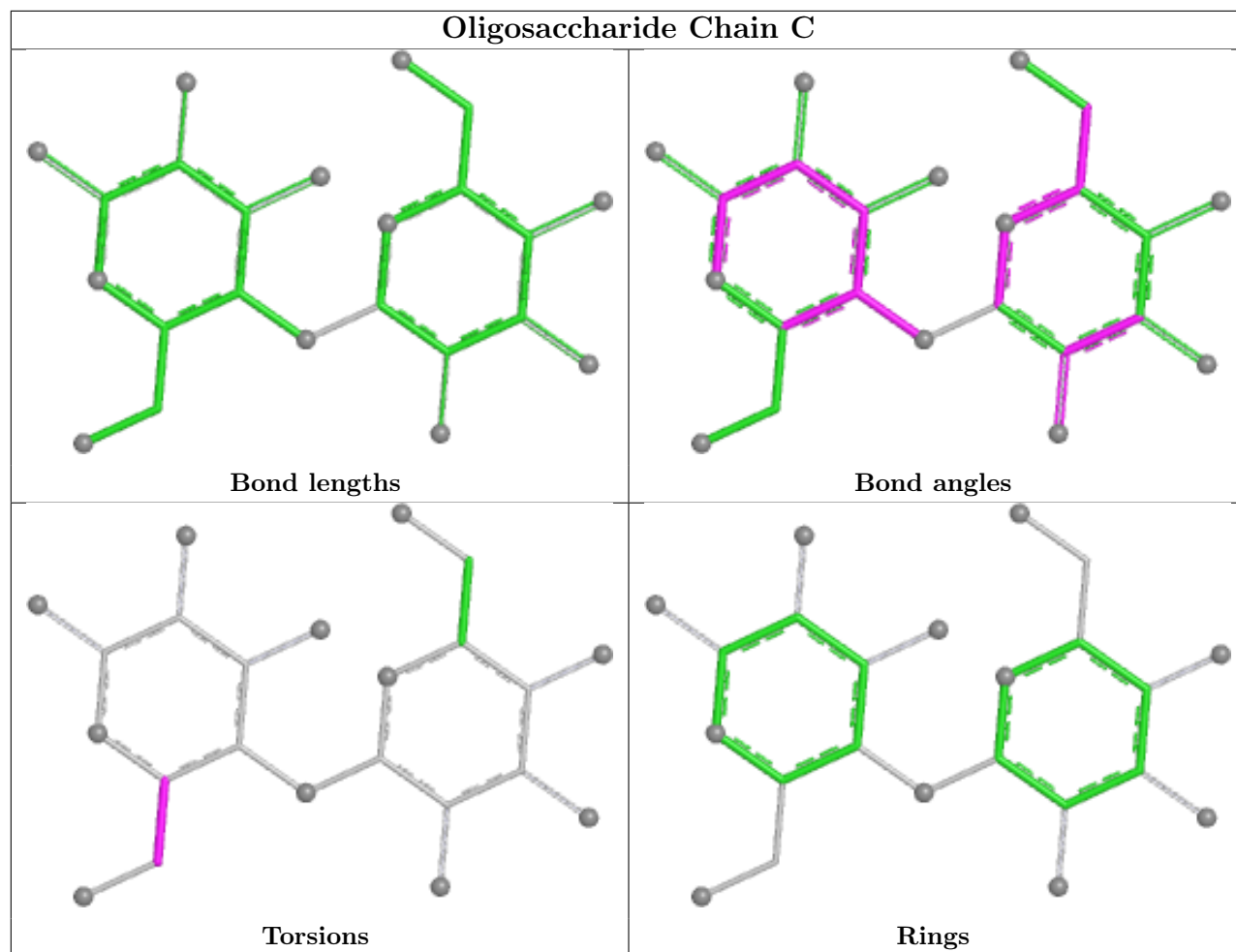
There are no ring outliers.

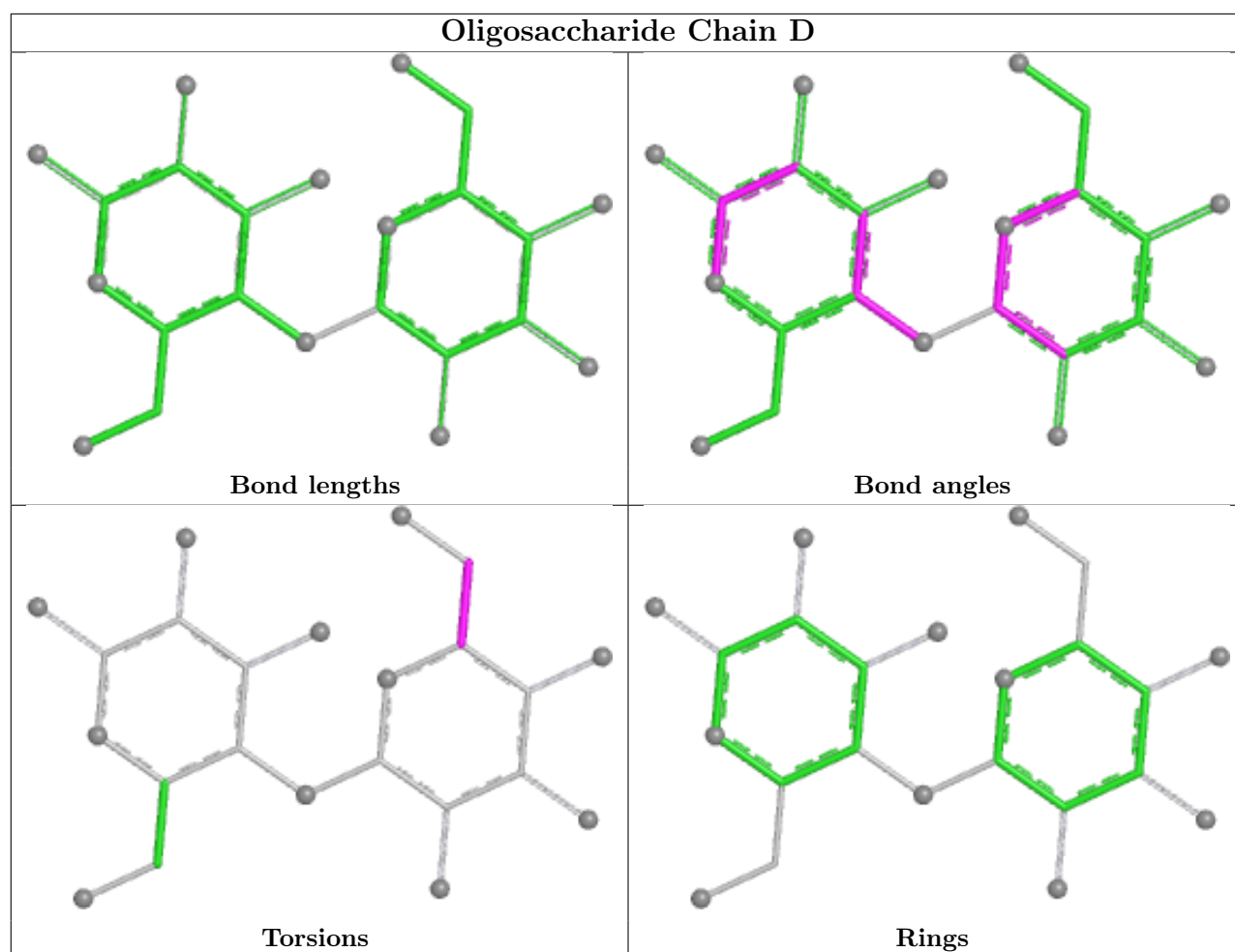
2 monomers are involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	D	1	GLC	1	0
2	D	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](#)

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 [i](#)

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%)	-0.72	2 (0%) 90 87	9, 22, 43, 81	0

All (2) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	600	LEU	2.1
1	A	148	GLN	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](#)

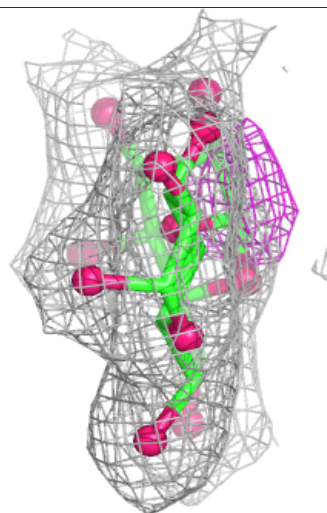
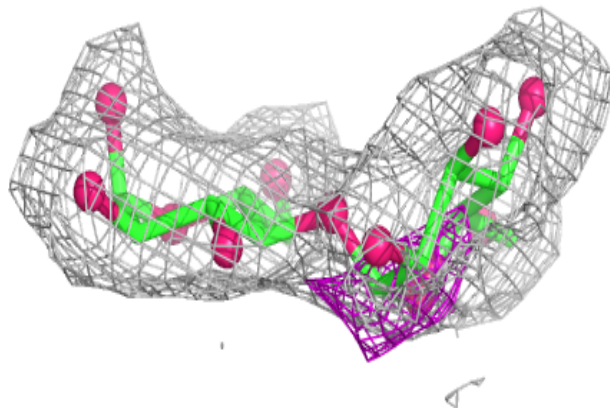
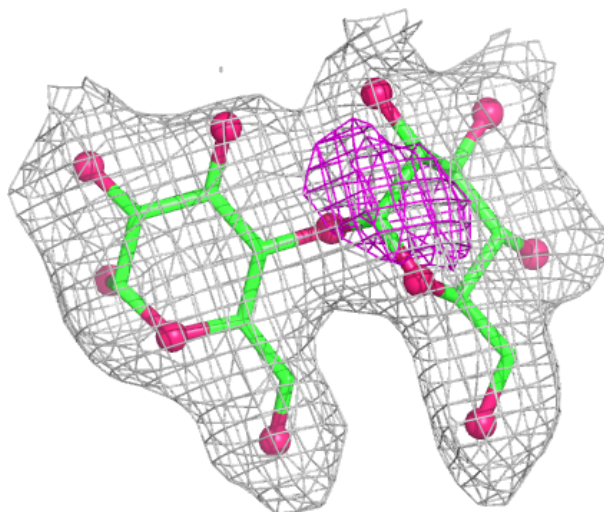
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled ‘Q< 0.9’ lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	GLC	D	2	11/12	0.69	0.20	98,98,100,100	0
2	GLC	D	1	12/12	0.72	0.17	91,93,96,99	0
2	GLC	C	1	12/12	0.88	0.13	52,53,57,63	0
2	GLC	B	2	11/12	0.90	0.10	24,27,29,32	0
2	GLC	C	2	11/12	0.90	0.10	48,50,55,55	0
2	GLC	B	1	12/12	0.95	0.07	18,24,30,32	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.

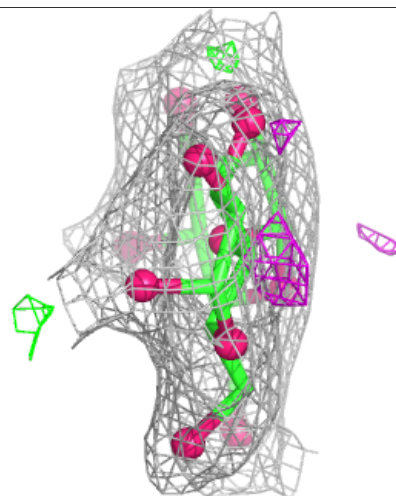
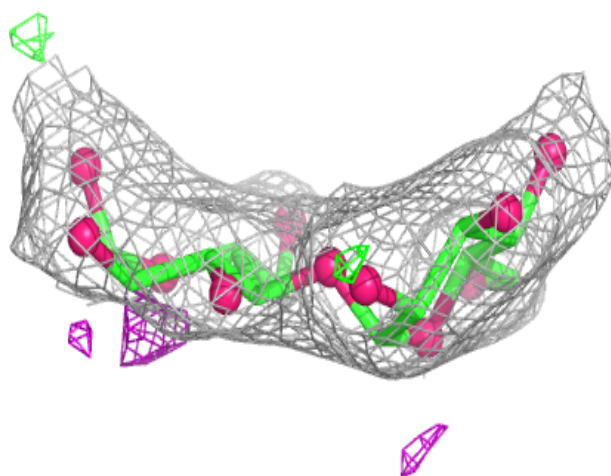
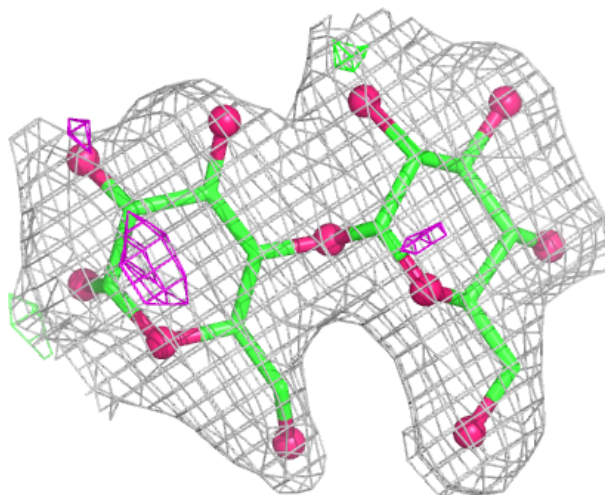
Electron density around Chain B:

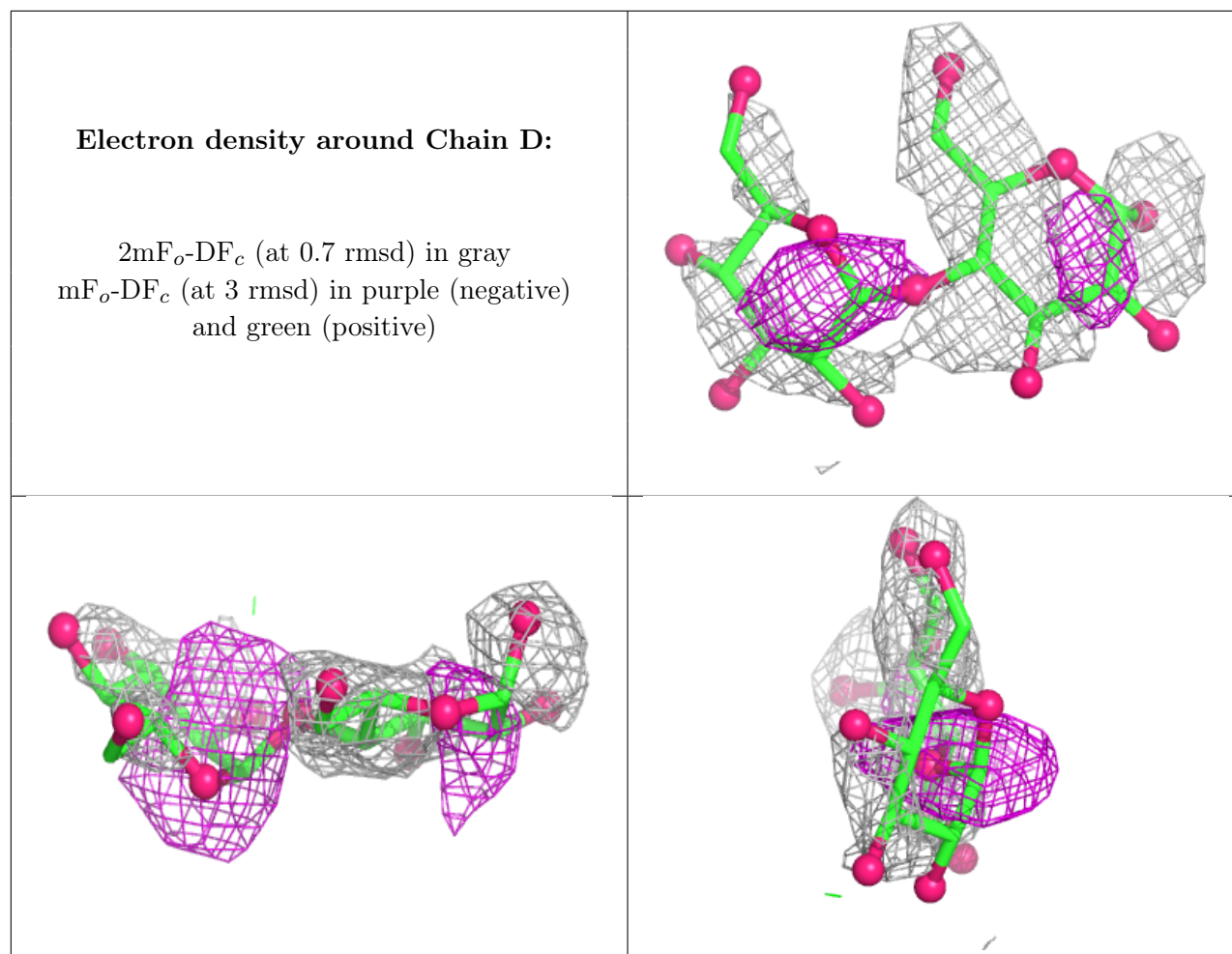
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around Chain C:

$2mF_o-DF_c$ (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
3	CA	A	691	1/1	1.00	0.07	22,22,22,22	0
3	CA	A	692	1/1	1.00	0.02	18,18,18,18	0

6.5 Other polymers [i](#)

There are no such residues in this entry.