



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 05:20 PM UTC

PDB ID : 3K8L / pdb_00003k8l
Title : Crystal structure of SusG-D498N mutant with maltoheptaose
Authors : Koropatkin, N.M.; Smith, T.J.
Deposited on : 2009-10-14
Resolution : 2.30 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

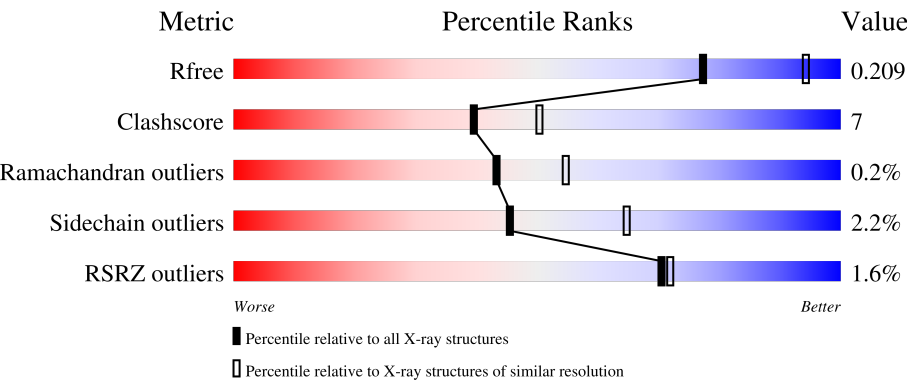
1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.30 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.



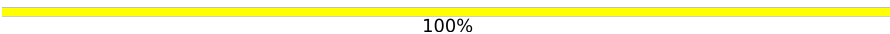
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R _{free}	180053	6319 (2.30-2.30)
Clashscore	190562	6919 (2.30-2.30)
Ramachandran outliers	187476	6854 (2.30-2.30)
Sidechain outliers	187428	6854 (2.30-2.30)
RSRZ outliers	180081	6325 (2.30-2.30)

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	669	% 82%14%..
1	B	669	2% 81%14%..
2	C	7	43%43%14%
3	D	5	20%60%20%
3	F	5	40%20%40%

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Mol	Chain	Length	Quality of chain
4	E	6	 33%50%17%
4	G	6	 100%
4	H	6	 33%50%17%

2 Entry composition [i](#)

There are 7 unique types of molecules in this entry. The entry contains 11264 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 Alpha-amylase, susG.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	649	Total	C	N	O	S	0	0	0
			5169	3288	836	1029	16			
1	B	646	Total	C	N	O	S	0	0	0
			5146	3277	832	1021	16			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	498	ASN	ASP	engineered mutation	UNP Q8A1G3
B	498	ASN	ASP	engineered mutation	UNP Q8A1G3

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



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace
2	C	7	Total	C	O	0	0	0
			78	42	36			

- Molecule 3 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
3	D	5	Total	C	O	1	0	0
			56	30	26			
3	F	5	Total	C	O	0	0	0
			56	30	26			

- Molecule 4 is an oligosaccharide called alpha-D-glucopyranose-(1-4)-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
4	E	6	Total	C	O	0	0	0
			67	36	31			
4	G	6	Total	C	O	0	0	0
			67	36	31			
4	H	6	Total	C	O	0	0	0
			67	36	31			

- 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		
5	B	2	Total	Ca	0	0
			2	2		

- Molecule 6 is 1,2-ETHANEDIOL (CCD ID: EDO) (formula: C₂H₆O₂).



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

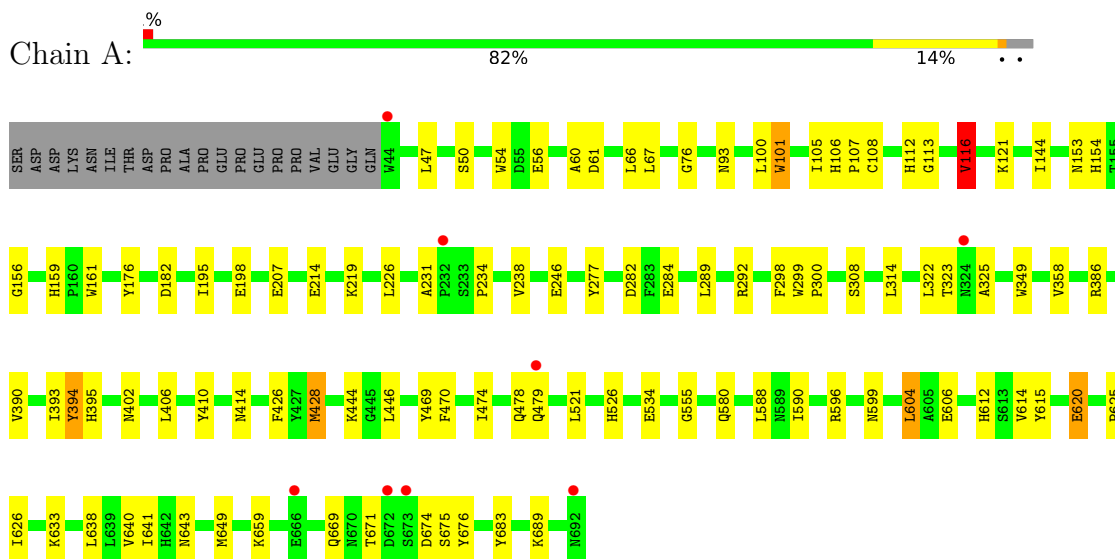
- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	342	Total	O	0	0
			342	342		
7	B	196	Total	O	0	0
			196	196		

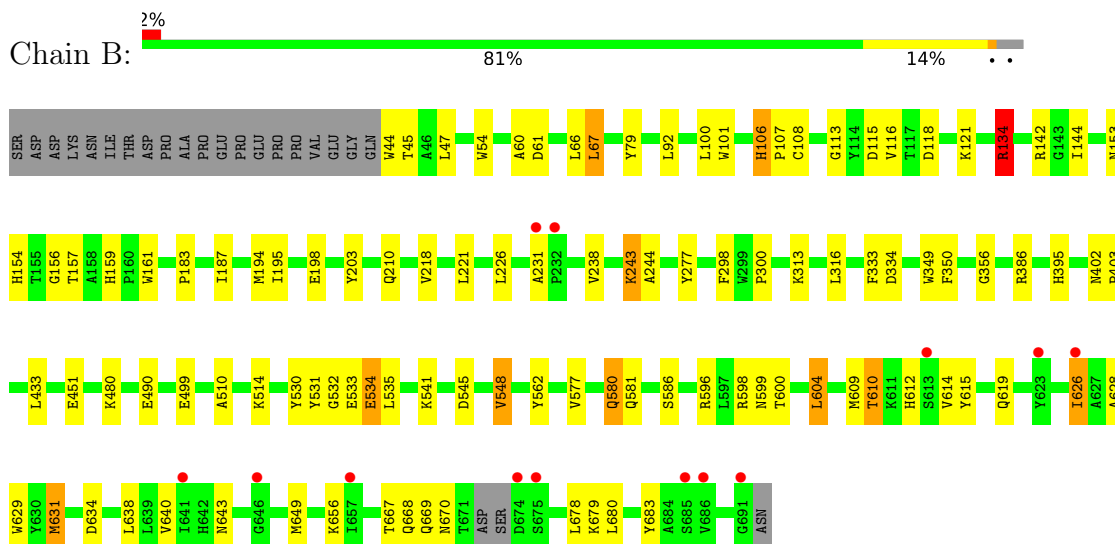
3 Residue-property plots

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.

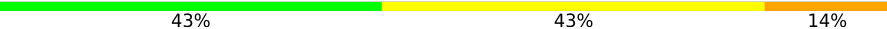
• Molecule 1: Alpha-amylase, susG

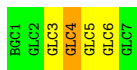


• Molecule 1: Alpha-amylase, susG

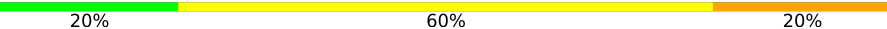


• Molecule 2: alpha-D-glucopyranose-(1-4)-alpha-D-glucopyranose-(1-4)-alpha-D-glucopyranose-(1-4)-alpha-D-glucopyranose-(1-4)-alpha-D-glucopyranose-(1-4)-beta-D-glucopyranose

Chain C:  43% 43% 14%

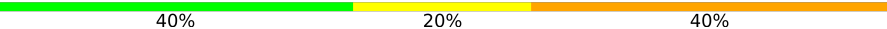


- Molecule 3: alpha-D-glucopyranose-(1-4)-alpha-D-glucopyranose-(1-4)-alpha-D-glucopyranose-(1-4)-alpha-D-glucopyranose-(1-4)-alpha-D-glucopyranose

Chain D:  20% 60% 20%



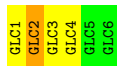
- Molecule 3: alpha-D-glucopyranose-(1-4)-alpha-D-glucopyranose-(1-4)-alpha-D-glucopyranose-(1-4)-alpha-D-glucopyranose-(1-4)-alpha-D-glucopyranose

Chain F:  40% 20% 40%

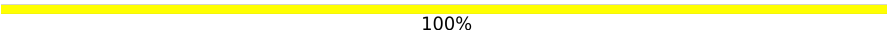


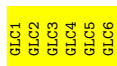
- Molecule 4: alpha-D-glucopyranose-(1-4)-alpha-D-glucopyranose-(1-4)-alpha-D-glucopyranose-(1-4)-alpha-D-glucopyranose-(1-4)-alpha-D-glucopyranose

Chain E:  33% 50% 17%

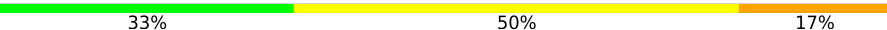


- Molecule 4: alpha-D-glucopyranose-(1-4)-alpha-D-glucopyranose-(1-4)-alpha-D-glucopyranose-(1-4)-alpha-D-glucopyranose-(1-4)-alpha-D-glucopyranose

Chain G:  100%



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

Chain H:  33% 50% 17%



4 Data and refinement statistics

Property	Value	Source
Space group	P 41	Depositor
Cell constants a, b, c, α , β , γ	128.04Å 128.04Å 130.48Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	50.00 – 2.30 50.00 – 2.30	Depositor EDS
% Data completeness (in resolution range)	94.4 (50.00-2.30) 94.4 (50.00-2.30)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.06	Depositor
$\langle I/\sigma(I) \rangle$ ¹	6.22 (at 2.20Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.184 , 0.216 0.179 , 0.209	Depositor DCC
R_{free} test set	9847 reflections (9.99%)	wwPDB-VP
Wilson B-factor (Å ²)	27.0	Xtriage
Anisotropy	0.508	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 31.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	0.018 for -h,-l,-k 0.005 for -h,l,k 0.006 for l,-k,h 0.019 for -l,-k,-h 0.032 for h,-k,-l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	11264	wwPDB-VP
Average B, all atoms (Å ²)	33.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.55% 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: GLC, BGC, EDO, CA

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	0.40	1/5308 (0.0%)	0.88	20/7214 (0.3%)
1	B	0.37	1/5284 (0.0%)	0.82	9/7181 (0.1%)
All	All	0.38	2/10592 (0.0%)	0.85	29/14395 (0.2%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	428	MET	SD-CE	-5.11	1.66	1.79
1	B	631	MET	SD-CE	-5.00	1.67	1.79

All (29) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	548	VAL	N-CA-C	-6.47	106.43	112.83
1	A	626	ILE	N-CA-C	6.20	117.72	108.54
1	B	626	ILE	N-CA-C	6.09	117.35	108.58
1	B	156	GLY	N-CA-C	-5.98	104.89	112.54
1	A	349	TRP	N-CA-C	-5.90	105.91	113.23
1	A	121	LYS	N-CA-C	5.81	117.44	109.18
1	A	246	GLU	N-CA-C	-5.78	101.93	110.48
1	A	106	HIS	N-CA-C	5.75	118.07	110.36
1	A	394	TYR	N-CA-C	-5.69	100.39	109.96
1	A	325	ALA	N-CA-C	5.66	118.22	110.24
1	B	106	HIS	N-CA-C	5.65	118.32	110.07
1	A	116	VAL	N-CA-C	5.62	117.28	109.80
1	A	689	LYS	N-CA-C	-5.56	100.17	109.24
1	A	393	ILE	N-CA-C	-5.55	104.92	110.30
1	A	358	VAL	N-CA-C	5.54	116.93	110.62
1	A	386	ARG	N-CA-C	-5.51	99.54	108.52
1	B	386	ARG	N-CA-C	-5.43	99.88	108.73

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	76	GLY	N-CA-C	5.35	122.58	114.61
1	A	282	ASP	N-CA-C	-5.35	99.25	108.56
1	B	121	LYS	N-CA-C	5.29	117.31	109.69
1	B	349	TRP	N-CA-C	-5.28	106.69	113.23
1	A	107	PRO	N-CA-C	-5.27	102.92	111.14
1	B	107	PRO	N-CA-C	-5.26	102.93	111.14
1	A	156	GLY	N-CA-C	-5.24	105.58	112.82
1	A	555	GLY	N-CA-C	-5.19	108.07	115.64
1	A	521	LEU	N-CA-C	5.09	119.34	113.18
1	B	134	ARG	N-CA-C	-5.02	105.88	111.36
1	A	182	ASP	CA-C-N	5.00	124.44	119.24
1	A	182	ASP	C-N-CA	5.00	124.44	119.24

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	5169	0	4872	61	0
1	B	5146	0	4856	74	0
2	C	78	0	66	1	0
3	D	56	0	48	1	0
3	F	56	0	48	1	0
4	E	67	0	57	2	0
4	G	67	0	57	0	0
4	H	67	0	57	2	0
5	A	2	0	0	0	0
5	B	2	0	0	0	0
6	A	12	0	18	2	0
6	B	4	0	6	1	0
7	A	342	0	0	3	0
7	B	196	0	0	2	0
All	All	11264	0	10085	139	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 7.

All (139) 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:394:TYR:H	1:A:402:ASN:HD21	1.04	0.93
1:B:159:HIS:HD2	1:B:161:TRP:H	1.16	0.92
1:A:159:HIS:HD2	1:A:161:TRP:H	1.14	0.92
1:A:153:ASN:HD22	1:A:154:HIS:HD2	1.16	0.88
1:B:153:ASN:HD22	1:B:154:HIS:HD2	1.19	0.88
1:B:609:MET:HB2	1:B:631:MET:HE1	1.54	0.87
1:B:313:LYS:HG2	1:B:334:ASP:HA	1.59	0.85
1:A:612:HIS:HD2	1:A:615:TYR:H	1.27	0.82
1:B:596:ARG:O	1:B:600:THR:HG23	1.80	0.82
1:B:243:LYS:HE2	1:B:244:ALA:H	1.47	0.79
1:A:604:LEU:HD13	1:A:638:LEU:HD12	1.65	0.79
1:A:238:VAL:HG21	1:A:314:LEU:HD21	1.65	0.79
1:B:54:TRP:CD1	1:B:600:THR:HG22	2.17	0.78
1:A:159:HIS:CD2	1:A:161:TRP:H	2.01	0.78
1:B:604:LEU:HD13	1:B:638:LEU:HD12	1.71	0.73
1:A:153:ASN:HD22	1:A:154:HIS:CD2	2.06	0.70
1:B:153:ASN:HD22	1:B:154:HIS:CD2	2.07	0.70
1:B:60:ALA:H	1:B:599:ASN:HD22	1.38	0.70
1:A:61:ASP:OD2	1:A:526:HIS:HD2	1.77	0.68
1:A:414:ASN:ND2	1:A:426:PHE:H	1.92	0.67
1:B:183:PRO:O	1:B:187:ILE:HG12	1.95	0.66
1:B:67:LEU:HD22	1:B:531:TYR:HE2	1.59	0.66
1:A:238:VAL:CG2	1:A:314:LEU:HD21	2.26	0.65
1:B:243:LYS:HE2	1:B:244:ALA:N	2.11	0.65
1:A:198:GLU:OE1	1:A:395:HIS:HD2	1.80	0.65
4:E:2:GLC:H62	4:E:3:GLC:H5	1.78	0.64
1:B:153:ASN:ND2	1:B:154:HIS:HD2	1.93	0.64
1:B:159:HIS:CD2	1:B:161:TRP:H	2.07	0.63
1:A:60:ALA:H	1:A:599:ASN:HD22	1.43	0.63
1:A:410:TYR:HD2	1:A:428:MET:HE2	1.65	0.62
4:H:3:GLC:H61	4:H:4:GLC:H5	1.80	0.62
1:A:47:LEU:HD21	1:A:669:GLN:HB3	1.82	0.62
1:A:394:TYR:H	1:A:402:ASN:ND2	1.87	0.61
1:A:428:MET:HE3	1:A:446:LEU:CD2	2.29	0.61
1:B:612:HIS:CD2	1:B:614:VAL:H	2.18	0.61
1:A:219:LYS:HG2	1:A:284:GLU:HG3	1.81	0.61
1:B:67:LEU:HD22	1:B:531:TYR:CE2	2.36	0.60
1:B:194:MET:HE3	1:B:356:GLY:H	1.66	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:671:THR:HA	1:A:675:SER:O	2.03	0.59
1:B:577:VAL:O	1:B:581:GLN:HG3	2.03	0.59
1:B:629:TRP:CD1	1:B:631:MET:HE3	2.38	0.58
1:B:612:HIS:HD2	1:B:615:TYR:H	1.50	0.58
1:A:428:MET:HE3	1:A:446:LEU:HD22	1.86	0.58
1:B:433:LEU:HD23	4:H:2:GLC:O4	2.03	0.57
1:B:298:PHE:O	1:B:300:PRO:HD3	2.04	0.57
1:B:60:ALA:H	1:B:599:ASN:ND2	2.03	0.57
1:B:198:GLU:OE1	1:B:395:HIS:HD2	1.89	0.56
1:B:598:ARG:NH1	7:B:875:HOH:O	2.26	0.55
1:A:410:TYR:HB2	1:A:428:MET:HE2	1.89	0.55
1:B:238:VAL:HG13	1:B:316:LEU:HD23	1.88	0.54
1:B:221:LEU:HD11	1:B:244:ALA:HB2	1.89	0.54
1:A:410:TYR:CD2	1:A:428:MET:HE2	2.40	0.54
1:A:50:SER:O	1:A:659:LYS:HE2	2.08	0.53
1:B:66:LEU:C	1:B:66:LEU:HD12	2.33	0.53
1:B:226:LEU:HB2	1:B:277:TYR:HB2	1.89	0.53
1:B:499:GLU:HA	1:B:541:LYS:HD2	1.90	0.53
1:A:625:PRO:HD2	1:A:649:MET:HE1	1.91	0.53
1:B:100:LEU:HD13	1:B:144:ILE:CG2	2.39	0.53
1:A:596:ARG:HD3	7:A:967:HOH:O	2.08	0.53
3:D:3:GLC:H62	3:D:4:GLC:H5	1.90	0.53
1:B:194:MET:HE3	1:B:356:GLY:N	2.24	0.52
1:B:106:HIS:HB3	1:B:118:ASP:O	2.10	0.52
1:B:142:ARG:HH11	1:B:142:ARG:HG3	1.75	0.51
1:B:619:GLN:NE2	1:B:619:GLN:HA	2.27	0.50
1:B:609:MET:CB	1:B:631:MET:HE1	2.36	0.50
1:A:234:PRO:HB2	1:A:322:LEU:HD12	1.92	0.50
1:A:580:GLN:NE2	1:A:588:LEU:H	2.10	0.50
1:A:640:VAL:O	1:A:641:ILE:HD13	2.11	0.50
1:B:626:ILE:N	1:B:626:ILE:HD12	2.27	0.50
1:B:92:LEU:HD13	1:B:100:LEU:HD21	1.94	0.50
1:A:469:TYR:OH	1:A:620:GLU:HG3	2.12	0.49
1:A:54:TRP:CE2	1:A:56:GLU:HA	2.48	0.49
1:B:667:THR:HG22	1:B:680:LEU:HD23	1.94	0.48
1:B:108:CYS:SG	1:B:113:GLY:HA2	2.53	0.48
1:A:390:VAL:HB	1:A:406:LEU:HD11	1.95	0.48
1:B:609:MET:HB2	1:B:631:MET:CE	2.35	0.48
1:A:470:PHE:O	1:A:474:ILE:HG12	2.14	0.48
1:B:45:THR:HG22	1:B:669:GLN:O	2.13	0.48
1:B:218:VAL:HG21	1:B:333:PHE:CD1	2.49	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:612:HIS:CD2	1:A:614:VAL:H	2.31	0.47
1:B:79:TYR:CE1	1:B:562:TYR:HB2	2.49	0.47
1:A:444:LYS:HE3	1:B:634:ASP:OD2	2.14	0.47
1:B:612:HIS:HD2	1:B:615:TYR:N	2.12	0.47
1:A:176:TYR:OH	6:A:900:EDO:H11	2.15	0.47
1:A:112:HIS:HE2	6:A:930:EDO:H11	1.79	0.47
1:A:633:LYS:HG2	7:B:701:HOH:O	2.14	0.47
1:B:609:MET:HA	1:B:631:MET:HE2	1.96	0.47
1:B:101:TRP:C	1:B:101:TRP:CD1	2.92	0.46
1:B:643:ASN:O	1:B:683:TYR:HA	2.15	0.46
1:A:606:GLU:OE2	1:A:633:LYS:HE3	2.16	0.46
1:A:195:ILE:HG12	1:A:395:HIS:HB2	1.98	0.46
1:B:115:ASP:HB3	1:B:350:PHE:CE1	2.50	0.46
1:B:134:ARG:HH11	1:B:134:ARG:HB3	1.80	0.46
1:B:402:ASN:HB2	1:B:403:PRO:HD3	1.98	0.46
1:B:668:GLN:HB2	1:B:679:LYS:HB3	1.97	0.45
3:F:3:GLC:H62	3:F:4:GLC:H5	1.98	0.45
1:A:101:TRP:C	1:A:101:TRP:CD1	2.93	0.45
1:B:451:GLU:OE1	1:B:490:GLU:OE2	2.34	0.45
1:B:668:GLN:OE1	1:B:679:LYS:HD3	2.17	0.45
1:B:656:LYS:O	1:B:656:LYS:HG3	2.17	0.45
1:B:510:ALA:O	1:B:514:LYS:HG3	2.17	0.44
1:A:474:ILE:O	1:A:478:GLN:HG3	2.17	0.44
1:B:100:LEU:HD13	1:B:144:ILE:HG21	1.98	0.44
1:A:292:ARG:HD2	1:A:299:TRP:CH2	2.52	0.44
2:C:3:GLC:C6	2:C:4:GLC:H5	2.48	0.44
1:A:54:TRP:CD2	1:A:56:GLU:HA	2.52	0.44
1:B:44:TRP:CE3	1:B:670:ASN:HB2	2.52	0.44
1:A:116:VAL:HG13	7:A:753:HOH:O	2.17	0.44
1:B:580:GLN:NE2	1:B:586:SER:HB2	2.33	0.44
1:B:47:LEU:HD12	1:B:667:THR:HB	2.00	0.44
1:B:183:PRO:HB2	1:B:203:TYR:CE1	2.53	0.43
1:A:580:GLN:HE21	1:A:580:GLN:HB3	1.65	0.43
1:A:643:ASN:O	1:A:683:TYR:HA	2.17	0.43
1:B:203:TYR:OH	6:B:920:EDO:H21	2.18	0.43
1:A:108:CYS:SG	1:A:113:GLY:HA2	2.58	0.43
1:A:226:LEU:HB2	1:A:277:TYR:HB2	2.01	0.43
1:A:100:LEU:HD13	1:A:144:ILE:HG21	2.01	0.43
1:B:530:TYR:O	1:B:533:GLU:HG2	2.19	0.43
1:A:66:LEU:HD12	1:A:66:LEU:C	2.43	0.43
1:A:105:ILE:HG12	1:A:105:ILE:O	2.18	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:195:ILE:HG12	1:B:395:HIS:HB2	2.00	0.42
1:B:628:ALA:HA	1:B:640:VAL:O	2.19	0.42
1:B:535:LEU:HD11	1:B:577:VAL:HG23	2.02	0.42
1:B:545:ASP:O	1:B:548:VAL:HG22	2.20	0.42
1:A:308:SER:HA	1:A:323:THR:HG21	2.02	0.42
1:A:410:TYR:CB	1:A:428:MET:HE2	2.49	0.42
4:E:1:GLC:H61	4:E:2:GLC:H5	2.01	0.42
1:A:633:LYS:HD2	7:A:831:HOH:O	2.19	0.42
1:A:207:GLU:OE1	1:A:395:HIS:HE1	2.02	0.41
1:A:298:PHE:O	1:A:300:PRO:HD3	2.20	0.41
1:B:609:MET:HG2	1:B:610:THR:N	2.35	0.41
1:A:612:HIS:CD2	1:A:615:TYR:H	2.19	0.41
1:B:532:GLY:N	1:B:534:GLU:OE2	2.54	0.41
1:A:60:ALA:HB3	1:A:599:ASN:ND2	2.37	0.40
1:A:674:ASP:O	1:A:674:ASP:CG	2.63	0.40
1:B:54:TRP:HD1	1:B:600:THR:HG22	1.76	0.40
1:B:115:ASP:HB3	1:B:350:PHE:HE1	1.87	0.40
1:A:479:GLN:CD	1:B:610:THR:HG21	2.46	0.40
1:A:669:GLN:OE1	1:A:676:TYR:HE1	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	647/669 (97%)	621 (96%)	25 (4%)	1 (0%)	43	55
1	B	642/669 (96%)	603 (94%)	38 (6%)	1 (0%)	43	55
All	All	1289/1338 (96%)	1224 (95%)	63 (5%)	2 (0%)	43	55

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	231	ALA
1	B	231	ALA

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	550/568 (97%)	540 (98%)	10 (2%)	51	70
1	B	547/568 (96%)	533 (97%)	14 (3%)	40	59
All	All	1097/1136 (97%)	1073 (98%)	24 (2%)	45	65

All (24) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	67	LEU
1	A	93	ASN
1	A	101	TRP
1	A	116	VAL
1	A	214	GLU
1	A	289	LEU
1	A	534	GLU
1	A	590	ILE
1	A	604	LEU
1	A	620	GLU
1	B	61	ASP
1	B	67	LEU
1	B	116	VAL
1	B	134	ARG
1	B	157	THR
1	B	210	GLN
1	B	243	LYS
1	B	480	LYS
1	B	534	GLU
1	B	580	GLN
1	B	604	LEU
1	B	610	THR

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Mol	Chain	Res	Type
1	B	649	MET
1	B	678	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (43) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	93	ASN
1	A	154	HIS
1	A	159	HIS
1	A	197	GLN
1	A	210	GLN
1	A	230	ASN
1	A	274	ASN
1	A	324	ASN
1	A	337	GLN
1	A	342	HIS
1	A	395	HIS
1	A	402	ASN
1	A	414	ASN
1	A	464	ASN
1	A	479	GLN
1	A	526	HIS
1	A	543	ASN
1	A	561	ASN
1	A	580	GLN
1	A	589	ASN
1	A	599	ASN
1	A	608	ASN
1	A	612	HIS
1	A	668	GLN
1	B	87	GLN
1	B	94	GLN
1	B	154	HIS
1	B	159	HIS
1	B	197	GLN
1	B	317	ASN
1	B	330	ASN
1	B	395	HIS
1	B	419	GLN
1	B	422	HIS
1	B	543	ASN
1	B	561	ASN

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Mol	Chain	Res	Type
1	B	580	GLN
1	B	581	GLN
1	B	589	ASN
1	B	599	ASN
1	B	608	ASN
1	B	612	HIS
1	B	619	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 ⓘ

35 monosaccharides are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
2	BGC	C	1	2	12,12,12	0.92	0	17,17,17	0.81	0
2	GLC	C	2	2	11,11,12	0.90	0	15,15,17	0.80	0
2	GLC	C	3	2	11,11,12	1.14	0	15,15,17	0.65	0
2	GLC	C	4	2	11,11,12	1.18	1 (9%)	15,15,17	0.92	0
2	GLC	C	5	2	11,11,12	1.34	3 (27%)	15,15,17	0.69	0
2	GLC	C	6	2	11,11,12	1.47	2 (18%)	15,15,17	1.07	0
2	GLC	C	7	2	11,11,12	1.21	0	15,15,17	0.66	0
3	GLC	D	1	3	12,12,12	0.94	0	17,17,17	0.72	0
3	GLC	D	2	3	11,11,12	1.23	2 (18%)	15,15,17	0.66	0
3	GLC	D	3	3	11,11,12	1.13	0	15,15,17	0.69	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	GLC	D	4	3	11,11,12	1.10	1 (9%)	15,15,17	1.03	1 (6%)
3	GLC	D	5	3	11,11,12	1.43	2 (18%)	15,15,17	0.76	0
4	GLC	E	1	4	12,12,12	1.11	0	17,17,17	0.72	0
4	GLC	E	2	4	11,11,12	1.23	1 (9%)	15,15,17	0.64	0
4	GLC	E	3	4	11,11,12	1.16	0	15,15,17	0.73	0
4	GLC	E	4	4	11,11,12	1.15	1 (9%)	15,15,17	0.74	0
4	GLC	E	5	4	11,11,12	1.19	0	15,15,17	0.79	0
4	GLC	E	6	4	11,11,12	1.24	0	15,15,17	0.68	0
3	GLC	F	1	3	12,12,12	0.92	0	17,17,17	0.72	0
3	GLC	F	2	3	11,11,12	1.11	0	15,15,17	0.66	0
3	GLC	F	3	3	11,11,12	1.23	1 (9%)	15,15,17	0.76	0
3	GLC	F	4	3	11,11,12	1.27	2 (18%)	15,15,17	0.82	0
3	GLC	F	5	3	11,11,12	1.38	2 (18%)	15,15,17	0.68	0
4	GLC	G	1	4	12,12,12	1.50	3 (25%)	17,17,17	0.76	0
4	GLC	G	2	4	11,11,12	1.33	1 (9%)	15,15,17	0.68	0
4	GLC	G	3	4	11,11,12	1.39	2 (18%)	15,15,17	0.92	1 (6%)
4	GLC	G	4	4	11,11,12	1.44	3 (27%)	15,15,17	0.71	0
4	GLC	G	5	4	11,11,12	1.28	1 (9%)	15,15,17	0.76	0
4	GLC	G	6	4	11,11,12	1.40	1 (9%)	15,15,17	0.66	0
4	GLC	H	1	4	12,12,12	1.14	0	17,17,17	0.69	0
4	GLC	H	2	4	11,11,12	1.09	0	15,15,17	0.60	0
4	GLC	H	3	4	11,11,12	1.31	0	15,15,17	0.62	0
4	GLC	H	4	4	11,11,12	1.12	1 (9%)	15,15,17	0.91	0
4	GLC	H	5	4	11,11,12	1.36	2 (18%)	15,15,17	0.63	0
4	GLC	H	6	4	11,11,12	1.24	0	15,15,17	0.70	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
2	BGC	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	C	3	2	-	0/2/19/22	0/1/1/1
2	GLC	C	4	2	-	0/2/19/22	0/1/1/1
2	GLC	C	5	2	-	0/2/19/22	0/1/1/1
2	GLC	C	6	2	-	2/2/19/22	0/1/1/1
2	GLC	C	7	2	-	0/2/19/22	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	GLC	D	1	3	-	2/2/22/22	0/1/1/1
3	GLC	D	2	3	-	0/2/19/22	0/1/1/1
3	GLC	D	3	3	-	0/2/19/22	0/1/1/1
3	GLC	D	4	3	-	2/2/19/22	0/1/1/1
3	GLC	D	5	3	-	0/2/19/22	0/1/1/1
4	GLC	E	1	4	-	1/2/22/22	0/1/1/1
4	GLC	E	2	4	-	2/2/19/22	0/1/1/1
4	GLC	E	3	4	-	0/2/19/22	0/1/1/1
4	GLC	E	4	4	-	0/2/19/22	0/1/1/1
4	GLC	E	5	4	-	2/2/19/22	0/1/1/1
4	GLC	E	6	4	-	0/2/19/22	0/1/1/1
3	GLC	F	1	3	-	1/2/22/22	0/1/1/1
3	GLC	F	2	3	-	0/2/19/22	0/1/1/1
3	GLC	F	3	3	-	0/2/19/22	0/1/1/1
3	GLC	F	4	3	-	0/2/19/22	0/1/1/1
3	GLC	F	5	3	-	2/2/19/22	0/1/1/1
4	GLC	G	1	4	-	0/2/22/22	0/1/1/1
4	GLC	G	2	4	-	2/2/19/22	0/1/1/1
4	GLC	G	3	4	-	2/2/19/22	0/1/1/1
4	GLC	G	4	4	-	2/2/19/22	0/1/1/1
4	GLC	G	5	4	-	2/2/19/22	0/1/1/1
4	GLC	G	6	4	-	0/2/19/22	0/1/1/1
4	GLC	H	1	4	-	0/2/22/22	0/1/1/1
4	GLC	H	2	4	-	0/2/19/22	0/1/1/1
4	GLC	H	3	4	-	1/2/19/22	0/1/1/1
4	GLC	H	4	4	-	0/2/19/22	0/1/1/1
4	GLC	H	5	4	-	0/2/19/22	0/1/1/1
4	GLC	H	6	4	-	0/2/19/22	0/1/1/1

All (32) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	4	GLC	O5-C1	2.73	1.48	1.43
4	G	3	GLC	C2-C3	2.61	1.56	1.52
4	G	3	GLC	O5-C1	2.52	1.47	1.43
4	G	2	GLC	C1-C2	2.49	1.58	1.52
4	G	1	GLC	O4-C4	2.41	1.48	1.43
3	F	5	GLC	C2-C3	2.39	1.56	1.52
4	G	4	GLC	O5-C5	2.38	1.48	1.43
4	H	4	GLC	O5-C1	2.30	1.47	1.43

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	5	GLC	O5-C1	2.29	1.47	1.43
2	C	5	GLC	O5-C5	2.29	1.47	1.43
3	F	4	GLC	O5-C1	2.27	1.47	1.43
3	D	5	GLC	C2-C3	2.27	1.56	1.52
2	C	6	GLC	C4-C3	2.26	1.58	1.52
2	C	6	GLC	O4-C4	2.25	1.48	1.43
4	G	1	GLC	C4-C5	2.25	1.57	1.53
3	D	5	GLC	O5-C1	2.24	1.47	1.43
4	G	6	GLC	O5-C1	2.23	1.47	1.43
3	F	3	GLC	C2-C3	2.20	1.55	1.52
4	G	4	GLC	O5-C1	2.19	1.47	1.43
3	D	2	GLC	C2-C3	2.15	1.55	1.52
3	F	5	GLC	O5-C1	2.15	1.47	1.43
4	G	1	GLC	C1-C2	2.14	1.57	1.52
2	C	5	GLC	C4-C5	2.13	1.57	1.53
3	D	2	GLC	O5-C1	2.09	1.47	1.43
4	E	2	GLC	O5-C1	2.09	1.47	1.43
4	G	5	GLC	O5-C1	2.07	1.47	1.43
3	F	4	GLC	C2-C3	2.04	1.55	1.52
4	G	4	GLC	C2-C3	2.03	1.55	1.52
3	D	4	GLC	O5-C1	2.03	1.47	1.43
4	H	5	GLC	C4-C5	2.02	1.57	1.53
4	E	4	GLC	O5-C1	2.01	1.47	1.43
4	H	5	GLC	O5-C5	2.00	1.47	1.43

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	D	4	GLC	C1-C2-C3	-2.11	106.58	109.64
4	G	3	GLC	C1-C2-C3	-2.01	106.72	109.64

There are no chirality outliers.

All (25) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	G	4	GLC	O5-C5-C6-O6
3	F	5	GLC	O5-C5-C6-O6
2	C	1	BGC	O5-C5-C6-O6
4	E	2	GLC	O5-C5-C6-O6
4	G	2	GLC	O5-C5-C6-O6
2	C	1	BGC	C4-C5-C6-O6
4	E	2	GLC	C4-C5-C6-O6

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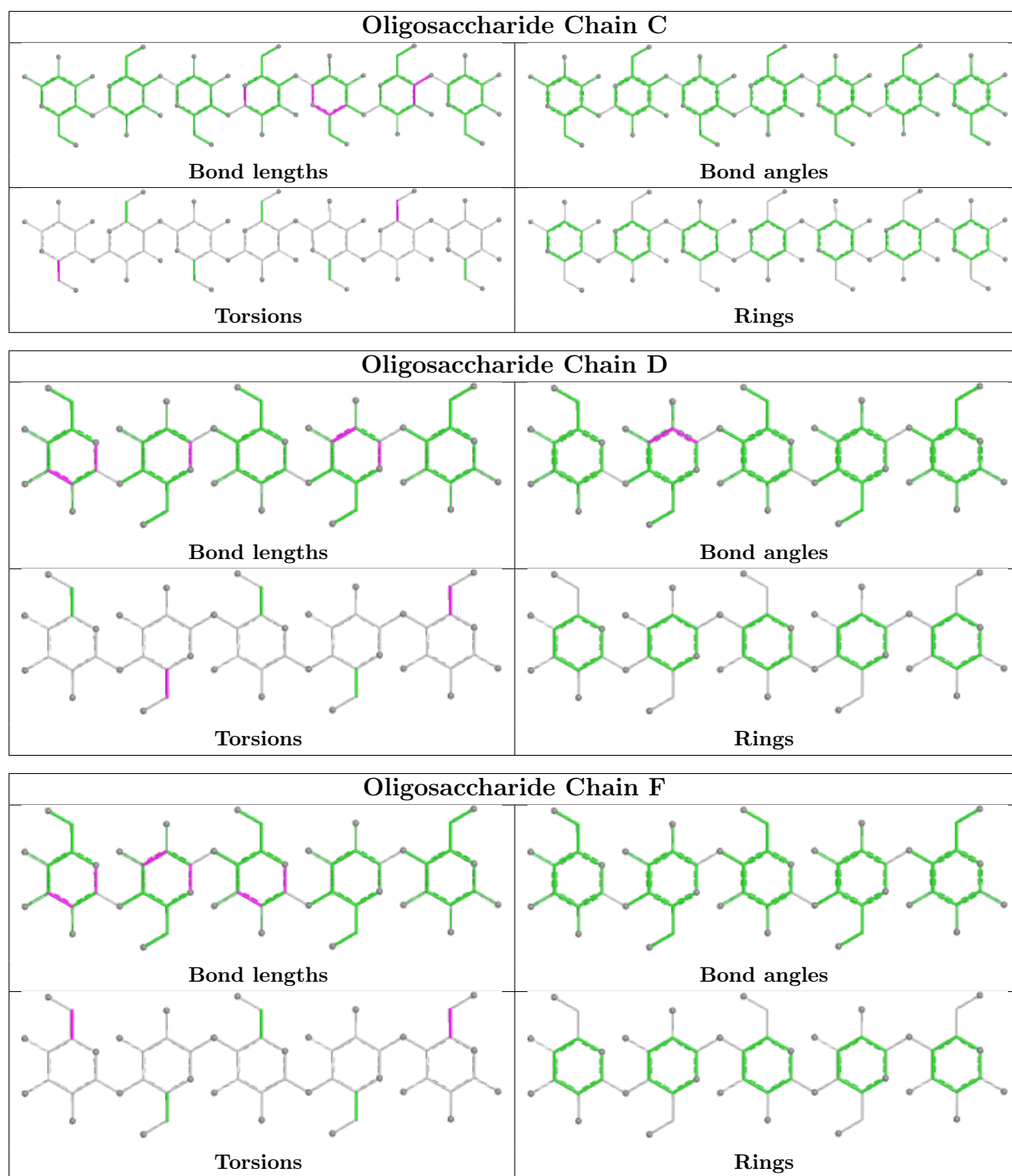
Mol	Chain	Res	Type	Atoms
4	G	5	GLC	O5-C5-C6-O6
3	F	5	GLC	C4-C5-C6-O6
4	G	4	GLC	C4-C5-C6-O6
4	G	5	GLC	C4-C5-C6-O6
3	D	1	GLC	O5-C5-C6-O6
3	D	4	GLC	O5-C5-C6-O6
3	D	4	GLC	C4-C5-C6-O6
4	G	2	GLC	C4-C5-C6-O6
3	D	1	GLC	C4-C5-C6-O6
4	E	5	GLC	O5-C5-C6-O6
4	G	3	GLC	C4-C5-C6-O6
4	E	1	GLC	O5-C5-C6-O6
3	F	1	GLC	O5-C5-C6-O6
2	C	6	GLC	C4-C5-C6-O6
4	G	3	GLC	O5-C5-C6-O6
2	C	6	GLC	O5-C5-C6-O6
4	H	3	GLC	O5-C5-C6-O6
4	E	5	GLC	C4-C5-C6-O6

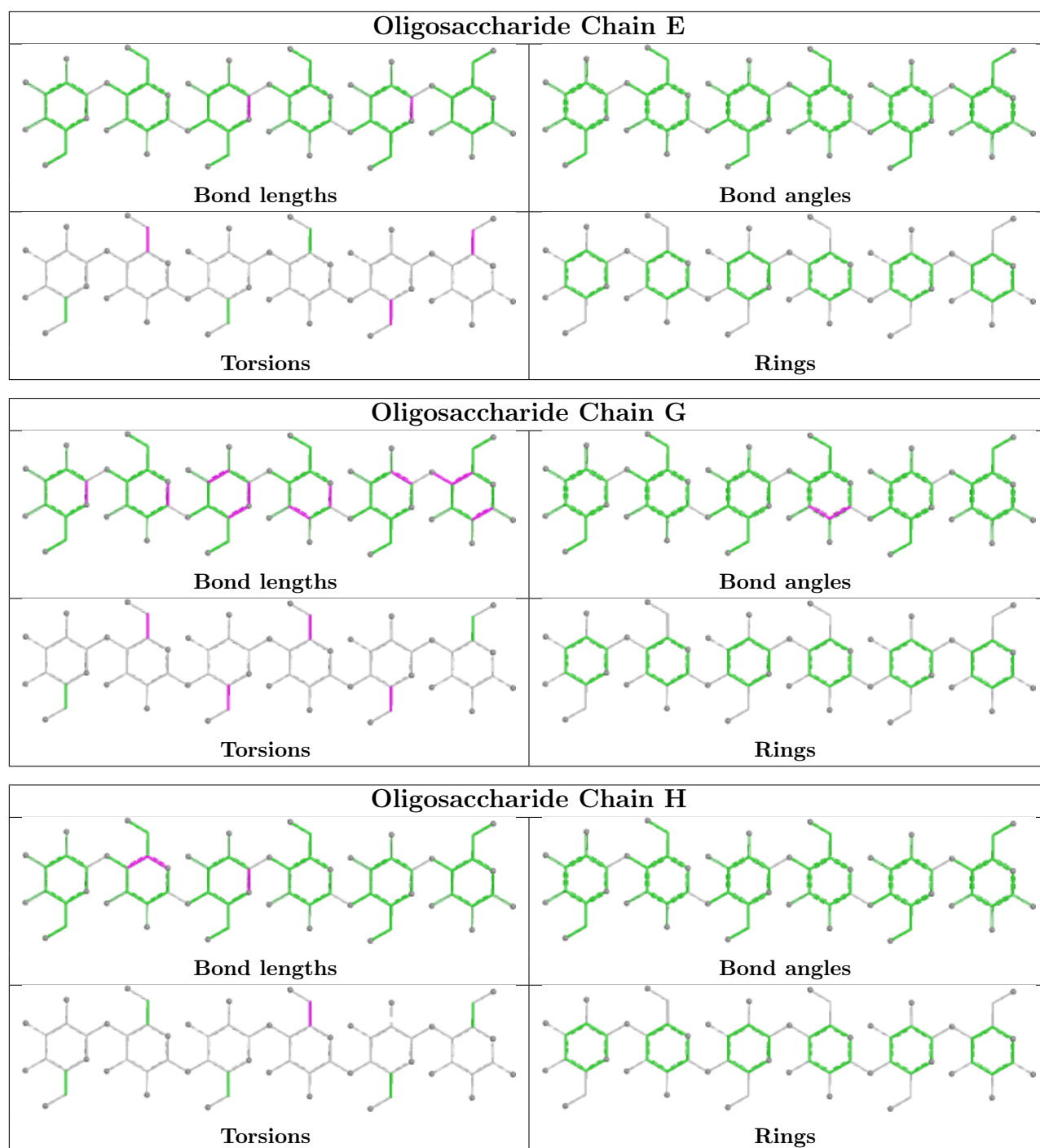
There are no ring outliers.

12 monomers are involved in 7 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	E	1	GLC	1	0
4	H	4	GLC	1	0
2	C	3	GLC	1	0
4	E	3	GLC	1	0
2	C	4	GLC	1	0
3	D	3	GLC	1	0
3	D	4	GLC	1	0
4	E	2	GLC	2	0
4	H	2	GLC	1	0
3	F	3	GLC	1	0
3	F	4	GLC	1	0
4	H	3	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 8 ligands modelled in this entry, 4 are monoatomic - leaving 4 for Mogul analysis.

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	$\# Z > 2$
6	EDO	A	930	-	3,3,3	0.71	0	2,2,2	0.39	0
6	EDO	A	910	-	3,3,3	0.62	0	2,2,2	0.42	0
6	EDO	A	900	-	3,3,3	0.70	0	2,2,2	0.39	0
6	EDO	B	920	-	3,3,3	0.74	0	2,2,2	0.36	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
6	EDO	A	930	-	-	0/1/1/1	-
6	EDO	A	910	-	-	0/1/1/1	-
6	EDO	A	900	-	-	0/1/1/1	-
6	EDO	B	920	-	-	1/1/1/1	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (1) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
6	B	920	EDO	O1-C1-C2-O2

There are no ring outliers.

3 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	A	930	EDO	1	0
6	A	900	EDO	1	0
6	B	920	EDO	1	0

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

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	649/669 (97%)	-0.37	8 (1%) 76 77	15, 26, 44, 67	0
1	B	646/669 (96%)	0.21	13 (2%) 65 66	20, 37, 54, 69	0
All	All	1295/1338 (96%)	-0.08	21 (1%) 70 72	15, 31, 52, 69	0

All (21) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	691	GLY	3.7
1	B	686	VAL	3.3
1	B	675	SER	3.1
1	B	674	ASP	3.0
1	A	692	ASN	2.9
1	A	673	SER	2.6
1	A	324	ASN	2.6
1	B	646	GLY	2.5
1	B	641	ILE	2.5
1	B	657	ILE	2.4
1	B	623	TYR	2.4
1	A	666	GLU	2.3
1	B	613	SER	2.2
1	A	232	PRO	2.1
1	A	479	GLN	2.1
1	B	685	SER	2.1
1	B	626	ILE	2.1
1	A	44	TRP	2.1
1	A	672	ASP	2.0
1	B	231	ALA	2.0
1	B	232	PRO	2.0

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

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

6.3 Carbohydrates ⓘ

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
4	GLC	G	6	11/12	0.50	0.18	72,75,76,76	0
3	GLC	F	5	11/12	0.60	0.20	65,68,70,71	0
4	GLC	G	2	11/12	0.64	0.19	60,66,68,68	0
4	GLC	H	6	11/12	0.64	0.23	59,62,65,66	0
3	GLC	D	5	11/12	0.65	0.22	61,65,66,66	0
4	GLC	H	1	12/12	0.68	0.20	61,70,71,72	0
4	GLC	G	1	12/12	0.68	0.20	66,69,70,71	0
4	GLC	G	3	11/12	0.76	0.15	56,57,58,58	0
4	GLC	E	6	11/12	0.77	0.14	61,64,66,68	0
2	BGC	C	1	12/12	0.81	0.17	48,58,60,61	0
3	GLC	D	4	11/12	0.81	0.15	41,46,50,55	1
4	GLC	G	5	11/12	0.81	0.12	59,62,64,68	0
4	GLC	E	5	11/12	0.82	0.13	44,49,56,57	0
2	GLC	C	6	11/12	0.83	0.14	40,43,51,52	0
2	GLC	C	7	11/12	0.83	0.17	47,51,52,53	0
4	GLC	E	2	11/12	0.85	0.12	40,50,54,55	0
3	GLC	F	1	12/12	0.87	0.12	31,43,45,47	0
4	GLC	E	1	12/12	0.87	0.13	53,55,57,59	0
4	GLC	H	4	11/12	0.90	0.10	32,37,39,41	0
4	GLC	H	2	11/12	0.91	0.11	47,52,56,56	0
3	GLC	D	1	12/12	0.91	0.10	38,42,44,46	0
4	GLC	E	3	11/12	0.91	0.10	30,32,36,38	0
3	GLC	F	4	11/12	0.92	0.10	41,48,52,58	0
4	GLC	H	5	11/12	0.92	0.10	40,41,47,53	0
4	GLC	G	4	11/12	0.92	0.09	51,54,55,56	0
4	GLC	H	3	11/12	0.93	0.09	38,39,42,43	0
2	GLC	C	5	11/12	0.93	0.09	22,23,29,34	0
2	GLC	C	4	11/12	0.94	0.07	21,23,25,26	0
4	GLC	E	4	11/12	0.94	0.08	31,34,36,38	0
2	GLC	C	3	11/12	0.95	0.07	24,27,29,35	0
2	GLC	C	2	11/12	0.95	0.08	35,40,43,45	0
3	GLC	D	2	11/12	0.95	0.07	30,32,33,35	0

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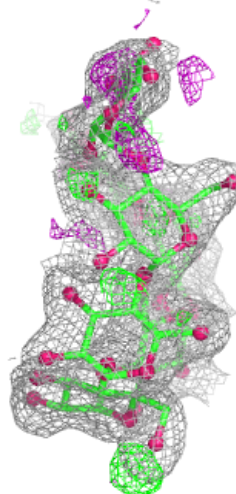
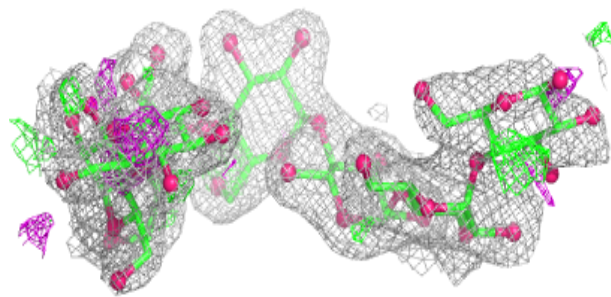
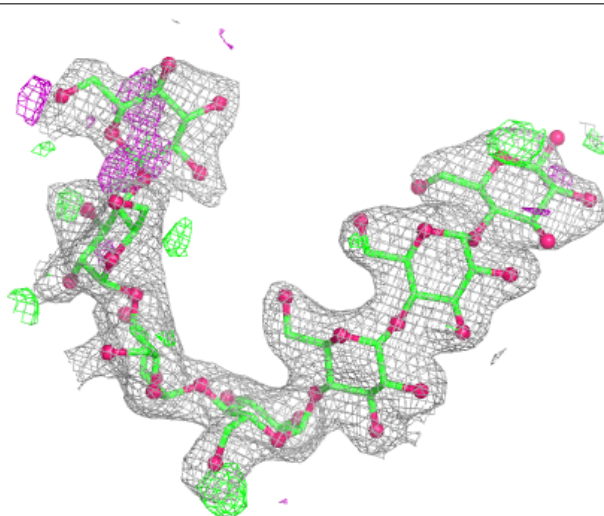
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	GLC	F	3	11/12	0.95	0.07	23,27,32,35	0
3	GLC	F	2	11/12	0.97	0.06	22,25,26,27	0
3	GLC	D	3	11/12	0.97	0.06	27,31,32,36	0

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.

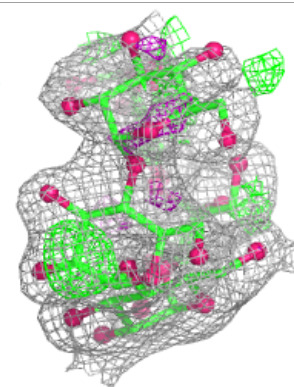
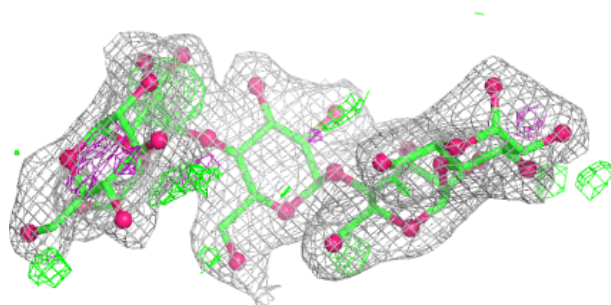
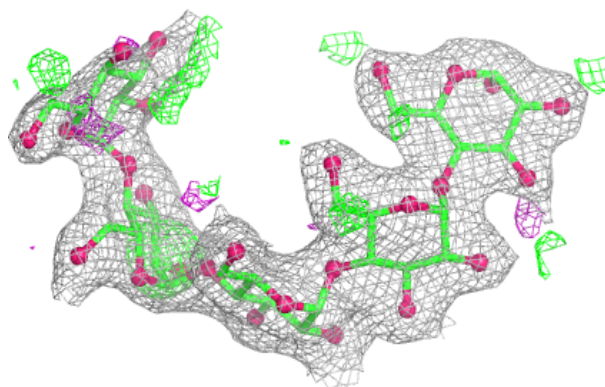
Electron density around Chain C:

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

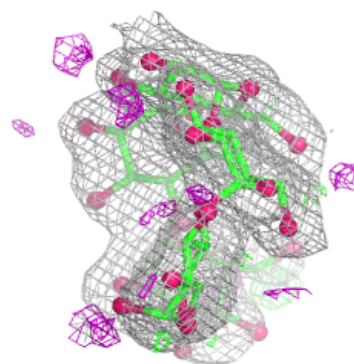
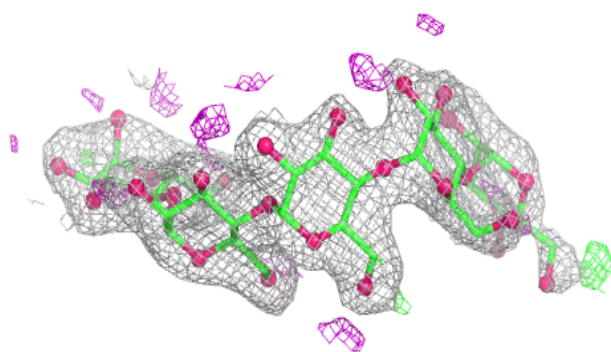
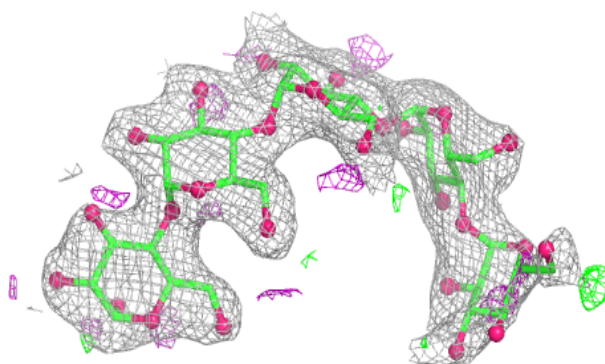


Electron density around Chain D:

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

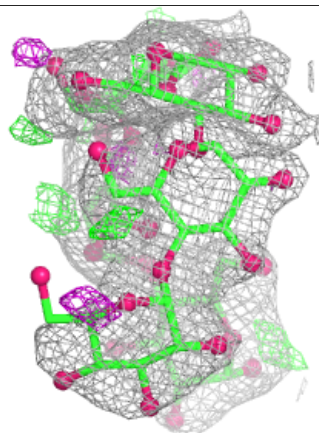
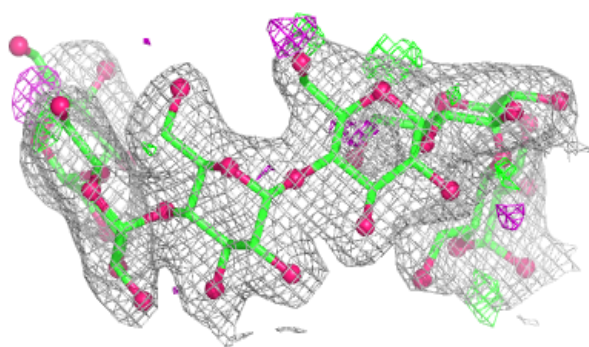
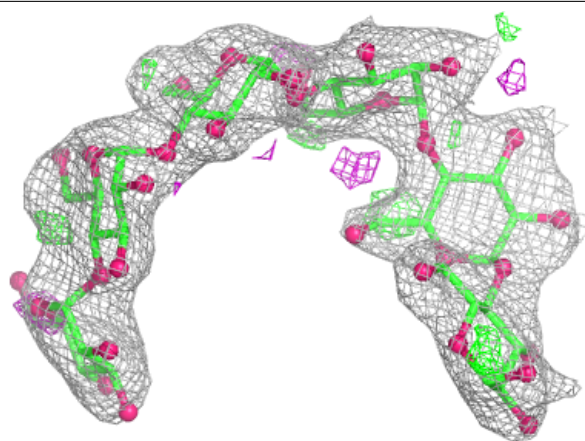
**Electron density around Chain F:**

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



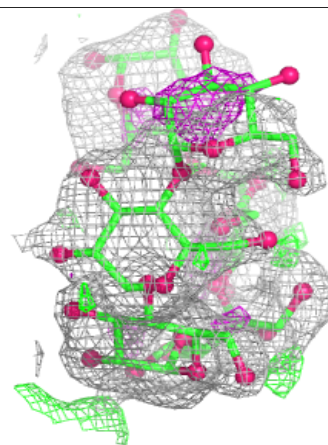
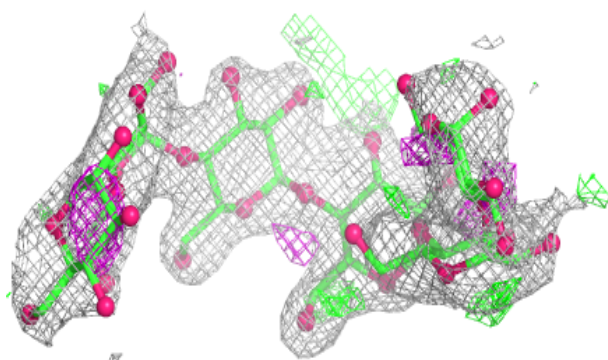
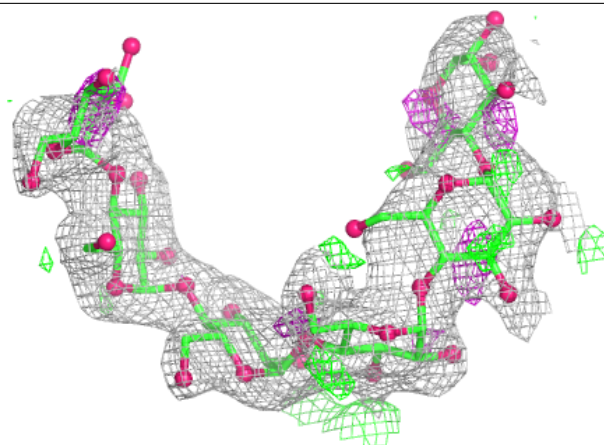
Electron density around Chain E:

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

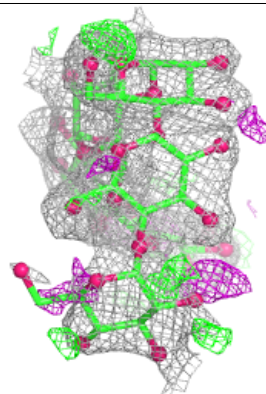
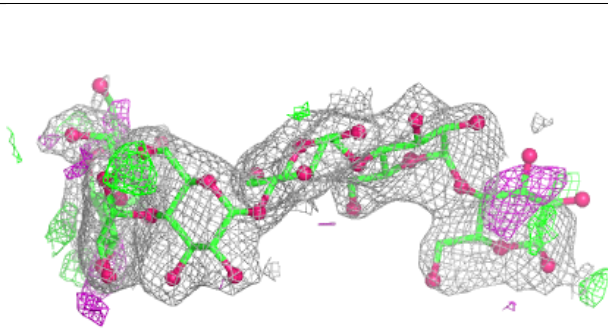
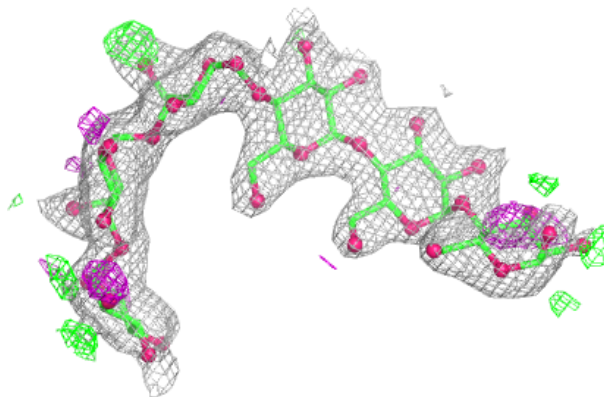


Electron density around Chain G:

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

**Electron density around Chain H:**

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



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
6	EDO	A	930	4/4	0.85	0.21	53,57,57,59	0
6	EDO	A	910	4/4	0.88	0.15	41,42,43,43	0
6	EDO	B	920	4/4	0.89	0.16	53,53,54,55	0
6	EDO	A	900	4/4	0.93	0.12	45,46,47,49	0
5	CA	A	700	1/1	0.98	0.04	31,31,31,31	0
5	CA	B	800	1/1	0.98	0.04	37,37,37,37	0
5	CA	A	710	1/1	0.99	0.03	16,16,16,16	0
5	CA	B	810	1/1	0.99	0.02	29,29,29,29	0

6.5 Other polymers [i](#)

There are no such residues in this entry.