



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 02:17 PM UTC

PDB ID : 6EPZ / pdb_00006epz
Title : Structure of the periplasmic binding protein MelB (Atu4661) in complex with melibiose from Agrobacterium fabrum C58
Authors : Vigouroux, A.; Morera, S.
Deposited on : 2017-10-12
Resolution : 1.80 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

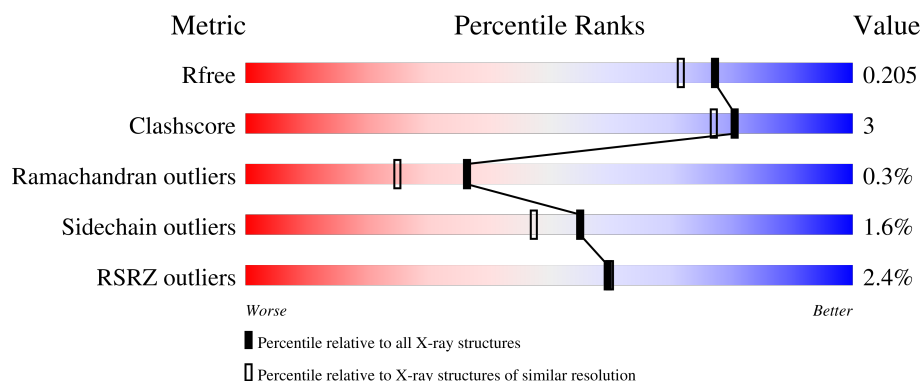
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

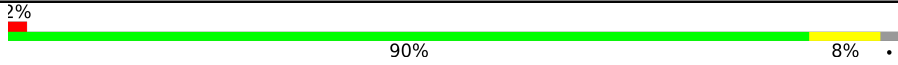
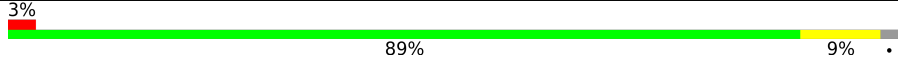
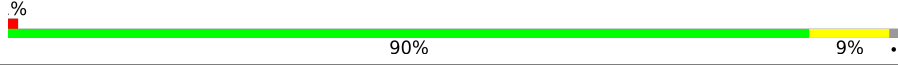
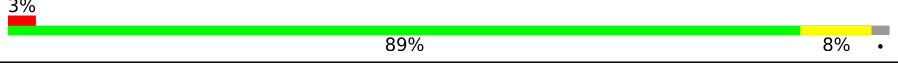
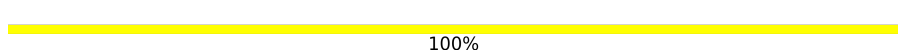
The reported resolution of this entry is 1.80 Å.

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	7662 (1.80-1.80)
Clashscore	190562	8479 (1.80-1.80)
Ramachandran outliers	187476	8391 (1.80-1.80)
Sidechain outliers	187428	8390 (1.80-1.80)
RSRZ outliers	180081	7663 (1.80-1.80)

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	683	
1	B	683	
1	C	683	
1	D	683	
2	E	2	

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Mol	Chain	Length	Quality of chain
2	F	2	 100%
2	G	2	 100%
2	H	2	 100%

2 Entry composition

There are 8 unique types of molecules in this entry. The entry contains 22767 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 Periplasmic alpha-galactoside-binding protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	672	Total	C	N	O	S	0	0	0
			5304	3386	902	1000	16			
1	D	671	Total	C	N	O	S	0	1	0
			5306	3386	903	1001	16			
1	C	673	Total	C	N	O	S	0	0	0
			5313	3390	904	1003	16			
1	B	671	Total	C	N	O	S	0	1	0
			5303	3384	901	1002	16			

There are 24 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	678	HIS	-	expression tag	UNP A0A083ZM57
A	679	HIS	-	expression tag	UNP A0A083ZM57
A	680	HIS	-	expression tag	UNP A0A083ZM57
A	681	HIS	-	expression tag	UNP A0A083ZM57
A	682	HIS	-	expression tag	UNP A0A083ZM57
A	683	HIS	-	expression tag	UNP A0A083ZM57
D	678	HIS	-	expression tag	UNP A0A083ZM57
D	679	HIS	-	expression tag	UNP A0A083ZM57
D	680	HIS	-	expression tag	UNP A0A083ZM57
D	681	HIS	-	expression tag	UNP A0A083ZM57
D	682	HIS	-	expression tag	UNP A0A083ZM57
D	683	HIS	-	expression tag	UNP A0A083ZM57
C	678	HIS	-	expression tag	UNP A0A083ZM57
C	679	HIS	-	expression tag	UNP A0A083ZM57
C	680	HIS	-	expression tag	UNP A0A083ZM57
C	681	HIS	-	expression tag	UNP A0A083ZM57
C	682	HIS	-	expression tag	UNP A0A083ZM57
C	683	HIS	-	expression tag	UNP A0A083ZM57
B	678	HIS	-	expression tag	UNP A0A083ZM57
B	679	HIS	-	expression tag	UNP A0A083ZM57
B	680	HIS	-	expression tag	UNP A0A083ZM57

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Chain	Residue	Modelled	Actual	Comment	Reference
B	681	HIS	-	expression tag	UNP A0A083ZM57
B	682	HIS	-	expression tag	UNP A0A083ZM57
B	683	HIS	-	expression tag	UNP A0A083ZM57

- Molecule 2 is an oligosaccharide called alpha-D-galactopyranose-(1-6)-alpha-D-glucopyranoside.



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace
2	E	2	Total	C	O	0	0	0
			23	12	11			
2	F	2	Total	C	O	0	0	0
			23	12	11			
2	G	2	Total	C	O	0	0	0
			23	12	11			
2	H	2	Total	C	O	0	0	0
			23	12	11			

- Molecule 3 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	Mg	0	0
			1	1		
3	D	1	Total	Mg	0	0
			1	1		
3	C	2	Total	Mg	0	0
			2	2		
3	B	1	Total	Mg	0	0
			1	1		

- Molecule 4 is 1,2-ETHANEDIOL (CCD ID: EDO) (formula: C₂H₆O₂).



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

- Molecule 5 is SODIUM ION (CCD ID: NA) (formula: Na).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	6	Total	Na	0	0
			6	6		
5	D	4	Total	Na	0	0
			4	4		
5	C	9	Total	Na	0	0
			9	9		

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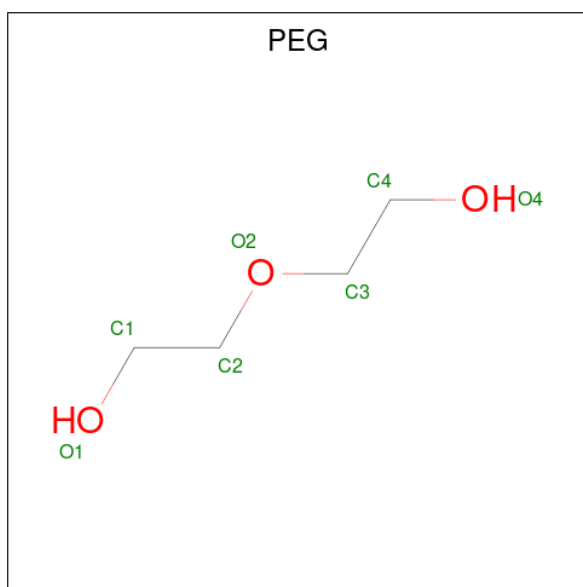
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	B	9	Total	Na	0	0
			9	9		

- Molecule 6 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	D	1	Total	Cl	0	0
			1	1		

- Molecule 7 is DI(HYDROXYETHYL)ETHER (CCD ID: PEG) (formula: C₄H₁₀O₃).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
7	D	1	Total	C	O	0	0
			7	4	3		

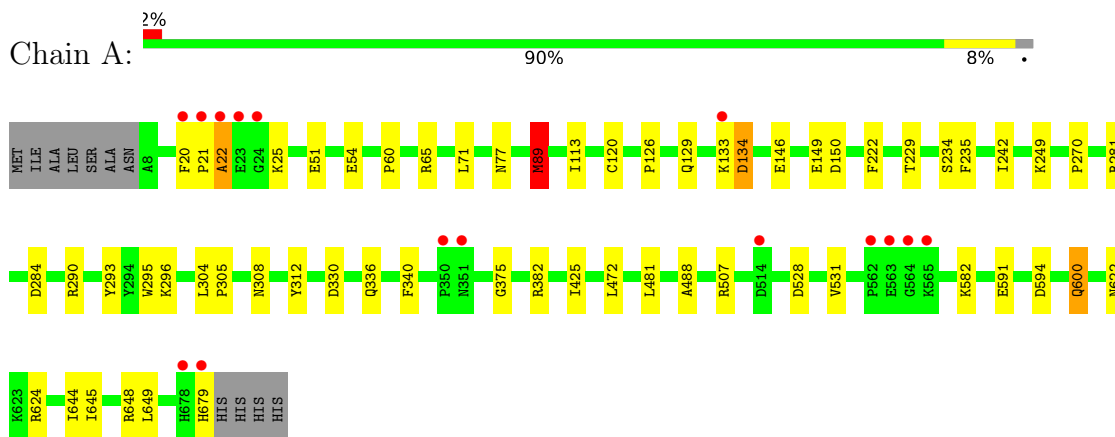
- Molecule 8 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	A	369	Total	O	0	0
			369	369		
8	D	332	Total	O	0	0
			332	332		
8	C	384	Total	O	0	0
			384	384		
8	B	287	Total	O	0	0
			287	287		

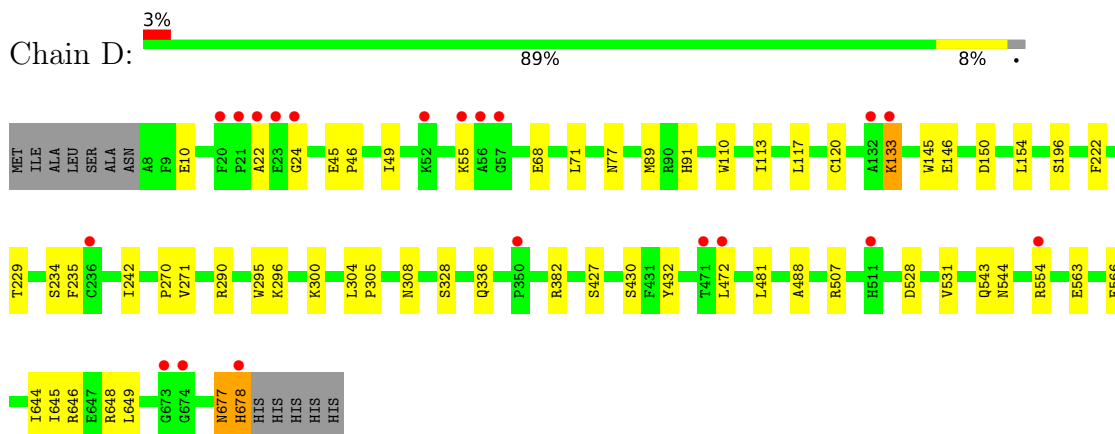
3 Residue-property plots [i](#)

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

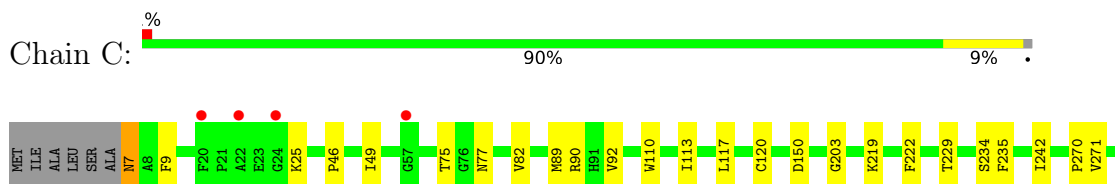
- Molecule 1: Periplasmic alpha-galactoside-binding protein

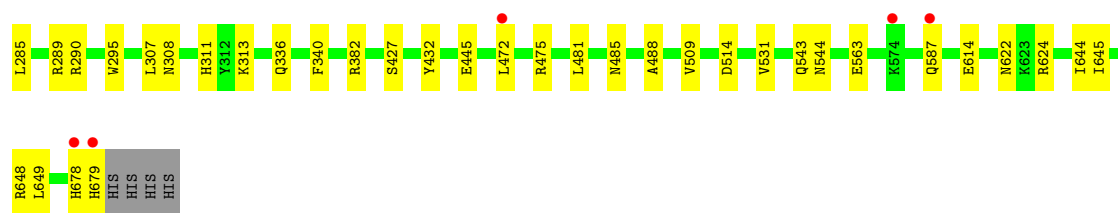


- Molecule 1: Periplasmic alpha-galactoside-binding protein

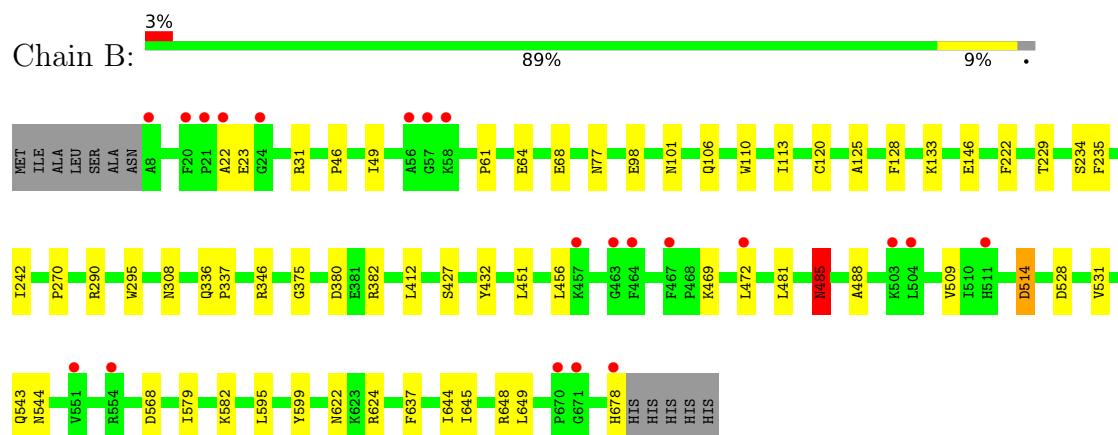


- Molecule 1: Periplasmic alpha-galactoside-binding protein





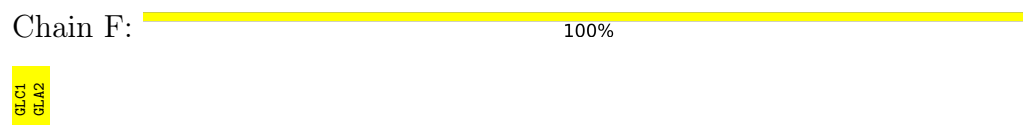
- Molecule 1: Periplasmic alpha-galactoside-binding protein



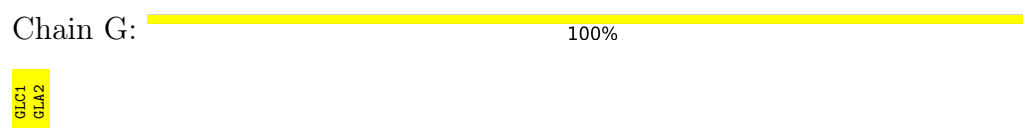
- Molecule 2: alpha-D-galactopyranose-(1-6)-alpha-D-glucopyranose



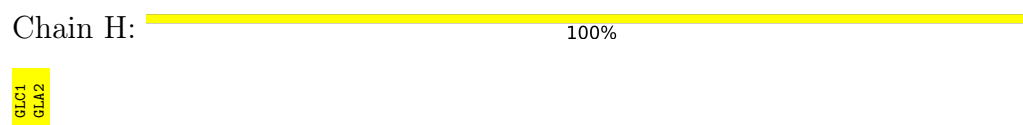
- Molecule 2: alpha-D-galactopyranose-(1-6)-alpha-D-glucopyranose



- Molecule 2: alpha-D-galactopyranose-(1-6)-alpha-D-glucopyranose



- Molecule 2: alpha-D-galactopyranose-(1-6)-alpha-D-glucopyranose



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	351.61Å 73.73Å 107.58Å 90.00° 105.38° 90.00°	Depositor
Resolution (Å)	48.85 – 1.80 48.85 – 1.80	Depositor EDS
% Data completeness (in resolution range)	99.9 (48.85-1.80) 99.9 (48.85-1.80)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.10	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.09 (at 1.79Å)	Xtriage
Refinement program	BUSTER 2.10.2	Depositor
R, R_{free}	0.177 , 0.195 0.186 , 0.205	Depositor DCC
R_{free} test set	12316 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	27.6	Xtriage
Anisotropy	0.400	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 39.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.28$	Xtriage
Estimated twinning fraction	0.026 for -h-2*k,l	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	22767	wwPDB-VP
Average B, all atoms (Å ²)	33.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 56.18 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 2.8422e-05. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

¹ Intensities estimated from amplitudes.

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

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: CL, PEG, MG, GLC, GLA, NA, EDO

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

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
1	A	0.85	2/5458 (0.0%)	1.08	5/7426 (0.1%)
1	B	0.81	1/5456 (0.0%)	1.09	9/7424 (0.1%)
1	C	0.84	1/5467 (0.0%)	1.08	7/7439 (0.1%)
1	D	0.82	1/5459 (0.0%)	1.09	8/7427 (0.1%)
All	All	0.83	5/21840 (0.0%)	1.08	29/29716 (0.1%)

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	89	MET	SD-CE	-10.90	1.52	1.79
1	C	222	PHE	CA-C	8.57	1.59	1.53
1	A	222	PHE	CA-C	6.21	1.57	1.53
1	D	222	PHE	CA-C	5.48	1.57	1.53
1	B	222	PHE	CA-C	5.41	1.57	1.53

All (29) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	514	ASP	CA-CB-CG	8.04	120.64	112.60
1	D	328	SER	CA-C-N	7.29	127.21	122.18
1	D	328	SER	C-N-CA	7.29	127.21	122.18
1	D	677	ASN	CA-C-N	7.20	134.65	121.70
1	D	677	ASN	C-N-CA	7.20	134.65	121.70
1	D	528	ASP	CA-CB-CG	6.38	118.98	112.60
1	B	485	ASN	CA-CB-CG	6.35	118.95	112.60
1	A	528	ASP	CA-CB-CG	6.18	118.78	112.60
1	B	528	ASP	CA-CB-CG	6.12	118.72	112.60
1	A	150	ASP	CA-CB-CG	5.51	118.11	112.60
1	B	128	PHE	N-CA-C	-5.42	105.77	112.38
1	C	514	ASP	CA-CB-CG	5.39	117.99	112.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	528	ASP	N-CA-C	-5.39	104.36	111.96
1	A	375	GLY	N-CA-C	5.38	122.86	115.27
1	A	340	PHE	CA-CB-CG	5.26	119.06	113.80
1	B	568	ASP	CA-CB-CG	5.25	117.85	112.60
1	C	563	GLU	CB-CG-CD	5.23	121.49	112.60
1	D	528	ASP	N-CA-C	-5.22	104.59	111.96
1	D	150	ASP	CA-CB-CG	5.20	117.80	112.60
1	B	380	ASP	CA-CB-CG	5.15	117.75	112.60
1	C	340	PHE	CA-CB-CG	5.13	118.93	113.80
1	C	485	ASN	CA-CB-CG	5.11	117.71	112.60
1	A	528	ASP	N-CA-C	-5.09	104.78	111.96
1	C	150	ASP	CA-CB-CG	5.08	117.68	112.60
1	B	375	GLY	N-CA-C	5.08	122.43	115.27
1	C	678	HIS	CA-C-N	5.05	130.80	121.70
1	C	678	HIS	C-N-CA	5.05	130.80	121.70
1	B	98	GLU	N-CA-C	-5.03	106.31	112.90
1	D	45	GLU	CB-CG-CD	5.02	121.13	112.60

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	5304	0	5119	30	0
1	B	5303	0	5117	27	0
1	C	5313	0	5125	27	0
1	D	5306	0	5124	25	0
2	E	23	0	21	0	0
2	F	23	0	21	0	0
2	G	23	0	21	0	0
2	H	23	0	21	0	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
3	C	2	0	0	0	0
3	D	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	A	16	0	24	0	0
4	B	8	0	12	1	0
4	C	4	0	6	0	0
4	D	8	0	12	1	0
5	A	6	0	0	0	0
5	B	9	0	0	0	0
5	C	9	0	0	0	0
5	D	4	0	0	0	0
6	D	1	0	0	0	0
7	D	7	0	10	0	0
8	A	369	0	0	0	0
8	B	287	0	0	0	0
8	C	384	0	0	1	0
8	D	332	0	0	0	0
All	All	22767	0	20633	107	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 3.

All (107) 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:425:ILE:H	1:A:600:GLN:HE22	1.18	0.91
1:C:92:VAL:HG22	1:C:313:LYS:HD2	1.67	0.77
1:A:77:ASN:HD22	1:A:648:ARG:HH12	1.36	0.74
1:C:77:ASN:HD22	1:C:648:ARG:HH12	1.37	0.71
1:B:77:ASN:HD22	1:B:648:ARG:HH12	1.37	0.71
1:D:77:ASN:HD22	1:D:648:ARG:HH12	1.40	0.70
1:C:120:CYS:HB3	1:C:235:PHE:O	1.94	0.68
1:D:120:CYS:HB3	1:D:235:PHE:O	1.94	0.68
1:A:120:CYS:HB3	1:A:235:PHE:O	1.95	0.67
1:B:120:CYS:HB3	1:B:235:PHE:O	1.95	0.67
1:B:61:PRO:HG2	1:B:64:GLU:HG3	1.82	0.62
1:A:89:MET:HE2	1:A:330:ASP:HB2	1.82	0.62
1:A:89:MET:CE	1:A:330:ASP:HB2	2.30	0.61
1:B:101:ASN:H	1:B:106:GLN:HE21	1.47	0.59
1:A:77:ASN:ND2	1:A:648:ARG:HH12	1.99	0.59
1:A:113:ILE:HG12	1:A:312:TYR:CD1	2.38	0.58
1:A:242:ILE:HG13	1:A:270:PRO:HG2	1.86	0.57
1:C:313:LYS:HD3	8:C:926:HOH:O	2.04	0.57
1:D:77:ASN:ND2	1:D:648:ARG:HH12	2.01	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:290:ARG:HH11	1:A:308:ASN:HD22	1.52	0.56
1:B:77:ASN:ND2	1:B:648:ARG:HH12	2.01	0.56
1:D:242:ILE:HG13	1:D:270:PRO:HG2	1.86	0.56
1:B:290:ARG:HH11	1:B:308:ASN:HD22	1.54	0.55
1:B:242:ILE:HG13	1:B:270:PRO:HG2	1.88	0.55
1:D:290:ARG:HH11	1:D:308:ASN:HD22	1.52	0.55
1:C:290:ARG:HH11	1:C:308:ASN:HD22	1.53	0.55
1:A:89:MET:CE	1:A:330:ASP:CB	2.85	0.54
1:C:7:ASN:HD22	1:C:9:PHE:H	1.54	0.54
1:C:77:ASN:ND2	1:C:648:ARG:HH12	2.03	0.53
1:B:31:ARG:HH22	4:B:702:EDO:H21	1.74	0.53
1:C:242:ILE:HG13	1:C:270:PRO:HG2	1.90	0.52
1:C:203:GLY:HA3	1:C:219:LYS:HE3	1.92	0.52
1:C:543:GLN:HE21	1:C:544:ASN:ND2	2.08	0.52
1:D:677:ASN:HB2	1:D:678:HIS:HB3	1.90	0.52
1:A:507:ARG:HD2	1:C:509:VAL:HG21	1.93	0.51
1:D:543:GLN:HE21	1:D:544:ASN:ND2	2.09	0.51
1:B:382:ARG:HA	1:B:472:LEU:HD11	1.93	0.51
1:D:481:LEU:HD11	1:D:531:VAL:HG23	1.93	0.51
1:B:543:GLN:HE21	1:B:544:ASN:ND2	2.09	0.50
1:A:425:ILE:N	1:A:600:GLN:HE22	1.98	0.49
1:C:472:LEU:O	1:C:475:ARG:HB2	2.12	0.49
1:A:20:PHE:C	1:A:22:ALA:H	2.21	0.49
1:A:229:THR:O	1:A:234:SER:HB2	2.13	0.49
1:A:600:GLN:HA	1:A:600:GLN:HE21	1.77	0.49
1:D:229:THR:O	1:D:234:SER:HB2	2.12	0.49
1:A:113:ILE:HG12	1:A:312:TYR:CG	2.48	0.48
1:B:110:TRP:O	1:B:113:ILE:HG22	2.13	0.48
1:A:65:ARG:HG2	1:A:293:TYR:CE1	2.49	0.48
1:C:46:PRO:HD2	1:C:49:ILE:HD12	1.95	0.48
1:C:229:THR:O	1:C:234:SER:HB2	2.14	0.48
1:C:89:MET:HB2	1:C:307:LEU:HD13	1.95	0.48
1:B:46:PRO:HD2	1:B:49:ILE:HD12	1.94	0.47
1:B:229:THR:O	1:B:234:SER:HB2	2.15	0.47
1:B:579:ILE:HG21	1:B:599:TYR:HB2	1.97	0.47
1:B:101:ASN:H	1:B:106:GLN:NE2	2.12	0.47
1:D:110:TRP:O	1:D:113:ILE:HG22	2.15	0.47
1:B:337:PRO:HB2	1:B:412:LEU:HB3	1.97	0.47
1:B:582:LYS:HD3	1:B:595:LEU:HD21	1.97	0.47
1:C:336:GLN:HE22	1:C:488:ALA:H	1.63	0.46
1:D:507:ARG:HD3	1:B:509:VAL:HG11	1.98	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:295:TRP:CD1	1:B:295:TRP:H	2.33	0.46
1:A:133:LYS:HA	1:A:134:ASP:HA	1.62	0.46
1:D:336:GLN:HE22	1:D:488:ALA:H	1.64	0.46
1:A:481:LEU:HD11	1:A:531:VAL:HG23	1.98	0.45
1:D:382:ARG:HA	1:D:472:LEU:HD11	1.97	0.45
1:C:481:LEU:HD11	1:C:531:VAL:HG23	1.98	0.45
1:A:382:ARG:HA	1:A:472:LEU:HD11	1.98	0.45
1:A:622:ASN:HD22	1:A:624:ARG:H	1.64	0.45
1:A:295:TRP:H	1:A:295:TRP:CD1	2.34	0.45
1:D:145:TRP:HB2	1:D:154:LEU:HD11	1.99	0.45
1:D:304:LEU:HB3	1:D:305:PRO:HA	1.97	0.45
1:A:71:LEU:HB3	1:A:296:LYS:HG2	1.99	0.45
1:A:304:LEU:HB3	1:A:305:PRO:HA	1.99	0.44
1:D:300:LYS:HB3	4:D:703:EDO:H21	1.98	0.44
1:C:117:LEU:HD22	1:C:271:VAL:HG11	1.99	0.44
1:B:427:SER:HA	1:B:432:TYR:CG	2.52	0.44
1:B:451:LEU:HD22	1:B:456:LEU:HD12	1.99	0.44
1:C:295:TRP:CD1	1:C:295:TRP:H	2.35	0.44
1:C:427:SER:HA	1:C:432:TYR:CG	2.52	0.44
1:C:622:ASN:HD22	1:C:624:ARG:H	1.66	0.44
1:B:481:LEU:HD11	1:B:531:VAL:HG23	1.99	0.44
1:C:110:TRP:O	1:C:113:ILE:HG22	2.18	0.43
1:D:295:TRP:CD1	1:D:295:TRP:H	2.35	0.43
1:D:196:SER:HB2	1:D:554:ARG:NH1	2.33	0.43
1:C:75:THR:HG22	1:C:82:VAL:HG13	2.00	0.43
1:A:336:GLN:HE22	1:A:488:ALA:H	1.65	0.43
1:A:644:ILE:HD12	1:A:649:LEU:HD11	1.99	0.43
1:D:46:PRO:HD2	1:D:49:ILE:HD12	2.01	0.43
1:C:382:ARG:HA	1:C:472:LEU:HD11	2.00	0.43
1:A:126:PRO:O	1:A:129:GLN:HG2	2.18	0.43
1:B:336:GLN:HE22	1:B:488:ALA:H	1.68	0.42
1:A:281:ARG:HD2	1:A:284:ASP:OD2	2.20	0.42
1:D:71:LEU:HB3	1:D:296:LYS:HG2	2.01	0.42
1:A:582:LYS:HE2	1:A:591:GLU:OE2	2.19	0.42
1:B:622:ASN:HD22	1:B:624:ARG:H	1.68	0.42
1:B:644:ILE:HD12	1:B:649:LEU:HD11	2.02	0.42
1:C:285:LEU:HD11	1:C:311:HIS:HB3	2.01	0.41
1:A:60:PRO:HD2	1:A:65:ARG:HD2	2.02	0.41
1:D:117:LEU:HD22	1:D:271:VAL:HG11	2.02	0.41
1:B:125:ALA:HA	1:B:637:PHE:HB2	2.03	0.41
1:D:91:HIS:HE2	1:D:646:ARG:HH11	1.67	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:133:LYS:HG2	1:D:430:SER:OG	2.21	0.41
1:C:289:ARG:HD2	1:C:289:ARG:HA	1.94	0.41
1:C:644:ILE:HD12	1:C:649:LEU:HD11	2.02	0.41
1:D:427:SER:HA	1:D:432:TYR:CG	2.56	0.41
1:D:644:ILE:HD12	1:D:649:LEU:HD11	2.03	0.40
1:B:485:ASN:ND2	1:B:514:ASP:HA	2.35	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	670/683 (98%)	646 (96%)	21 (3%)	3 (0%)	30	19
1	B	670/683 (98%)	646 (96%)	22 (3%)	2 (0%)	36	25
1	C	671/683 (98%)	650 (97%)	20 (3%)	1 (0%)	48	34
1	D	670/683 (98%)	645 (96%)	22 (3%)	3 (0%)	30	19
All	All	2681/2732 (98%)	2587 (96%)	85 (3%)	9 (0%)	36	25

All (9) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	22	ALA
1	D	22	ALA
1	D	24	GLY
1	B	22	ALA
1	A	21	PRO
1	A	645	ILE
1	C	645	ILE
1	B	645	ILE
1	D	645	ILE

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	558/568 (98%)	547 (98%)	11 (2%)	48	38
1	B	559/568 (98%)	551 (99%)	8 (1%)	59	52
1	C	560/568 (99%)	553 (99%)	7 (1%)	61	54
1	D	559/568 (98%)	550 (98%)	9 (2%)	55	47
All	All	2236/2272 (98%)	2201 (98%)	35 (2%)	55	47

All (35) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	25	LYS
1	A	51	GLU
1	A	54	GLU
1	A	89	MET
1	A	134	ASP
1	A	146	GLU
1	A	149	GLU
1	A	249	LYS
1	A	594	ASP
1	A	600	GLN
1	A	679	HIS
1	D	10	GLU
1	D	55	LYS
1	D	68	GLU
1	D	89	MET
1	D	133	LYS
1	D	146	GLU
1	D	563	GLU
1	D	566	GLU
1	D	678	HIS
1	C	7	ASN
1	C	25	LYS
1	C	90	ARG
1	C	445	GLU

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Mol	Chain	Res	Type
1	C	587	GLN
1	C	614	GLU
1	C	679	HIS
1	B	23	GLU
1	B	68	GLU
1	B	133	LYS
1	B	146	GLU
1	B	346	ARG
1	B	469	LYS
1	B	485	ASN
1	B	678	HIS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (61) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	77	ASN
1	A	246	GLN
1	A	253	ASN
1	A	302	GLN
1	A	308	ASN
1	A	336	GLN
1	A	339	ASN
1	A	466	ASN
1	A	516	ASN
1	A	543	GLN
1	A	544	ASN
1	A	547	GLN
1	A	598	GLN
1	A	600	GLN
1	A	622	ASN
1	A	677	ASN
1	D	77	ASN
1	D	302	GLN
1	D	308	ASN
1	D	323	GLN
1	D	336	GLN
1	D	351	ASN
1	D	368	GLN
1	D	500	GLN
1	D	516	ASN
1	D	526	GLN
1	D	544	ASN

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Mol	Chain	Res	Type
1	D	547	GLN
1	D	605	GLN
1	D	622	ASN
1	D	630	GLN
1	C	7	ASN
1	C	27	ASN
1	C	77	ASN
1	C	246	GLN
1	C	253	ASN
1	C	303	GLN
1	C	308	ASN
1	C	323	GLN
1	C	336	GLN
1	C	351	ASN
1	C	368	GLN
1	C	476	ASN
1	C	544	ASN
1	C	547	GLN
1	C	622	ASN
1	B	77	ASN
1	B	106	GLN
1	B	303	GLN
1	B	308	ASN
1	B	323	GLN
1	B	336	GLN
1	B	339	ASN
1	B	351	ASN
1	B	485	ASN
1	B	517	GLN
1	B	544	ASN
1	B	598	GLN
1	B	605	GLN
1	B	622	ASN
1	B	630	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 ⓘ

8 monosaccharides are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	GLC	E	1	2	12,12,12	1.47	2 (16%)	17,17,17	1.82	4 (23%)
2	GLA	E	2	2	11,11,12	1.26	2 (18%)	15,15,17	0.94	0
2	GLC	F	1	2	12,12,12	1.23	2 (16%)	17,17,17	1.20	1 (5%)
2	GLA	F	2	2	11,11,12	1.31	1 (9%)	15,15,17	1.06	1 (6%)
2	GLC	G	1	2	12,12,12	1.22	1 (8%)	17,17,17	1.59	3 (17%)
2	GLA	G	2	2	11,11,12	2.07	4 (36%)	15,15,17	1.30	1 (6%)
2	GLC	H	1	2	12,12,12	1.18	1 (8%)	17,17,17	1.76	3 (17%)
2	GLA	H	2	2	11,11,12	1.77	2 (18%)	15,15,17	1.03	1 (6%)

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	GLC	E	1	2	-	0/2/22/22	0/1/1/1
2	GLA	E	2	2	-	0/2/19/22	0/1/1/1
2	GLC	F	1	2	-	0/2/22/22	0/1/1/1
2	GLA	F	2	2	-	0/2/19/22	0/1/1/1
2	GLC	G	1	2	-	0/2/22/22	0/1/1/1
2	GLA	G	2	2	-	0/2/19/22	0/1/1/1
2	GLC	H	1	2	-	0/2/22/22	0/1/1/1
2	GLA	H	2	2	-	0/2/19/22	0/1/1/1

All (15) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	H	2	GLA	O5-C1	4.28	1.50	1.43
2	G	2	GLA	O5-C5	4.25	1.51	1.43

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	E	1	GLC	C1-C2	3.37	1.60	1.52
2	G	2	GLA	C2-C3	3.29	1.57	1.52
2	G	2	GLA	O5-C1	3.19	1.49	1.43
2	G	1	GLC	C1-C2	3.10	1.59	1.52
2	F	2	GLA	C2-C3	3.00	1.57	1.52
2	H	1	GLC	C1-C2	2.91	1.59	1.52
2	E	2	GLA	C1-C2	2.47	1.58	1.52
2	E	1	GLC	O5-C1	2.44	1.48	1.42
2	F	1	GLC	C1-C2	2.38	1.57	1.52
2	E	2	GLA	C2-C3	2.26	1.55	1.52
2	H	2	GLA	C2-C3	2.17	1.55	1.52
2	G	2	GLA	C4-C3	2.17	1.58	1.52
2	F	1	GLC	O5-C1	2.13	1.48	1.42

All (14) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	E	1	GLC	C1-O5-C5	4.34	122.05	113.65
2	H	1	GLC	C1-O5-C5	4.25	121.87	113.65
2	G	1	GLC	C1-O5-C5	3.81	121.02	113.65
2	G	2	GLA	O5-C5-C6	3.41	114.30	107.66
2	H	2	GLA	O5-C5-C6	3.07	113.64	107.66
2	F	1	GLC	C1-O5-C5	3.06	119.58	113.65
2	H	1	GLC	O5-C5-C4	2.92	114.96	109.70
2	H	1	GLC	O5-C5-C6	2.82	113.42	106.44
2	F	2	GLA	O5-C5-C6	2.78	113.07	107.66
2	E	1	GLC	O5-C5-C6	2.69	113.12	106.44
2	E	1	GLC	O5-C1-C2	2.64	114.95	110.30
2	G	1	GLC	O5-C5-C6	2.64	112.98	106.44
2	E	1	GLC	O5-C5-C4	2.64	114.45	109.70
2	G	1	GLC	O5-C1-C2	2.47	114.64	110.30

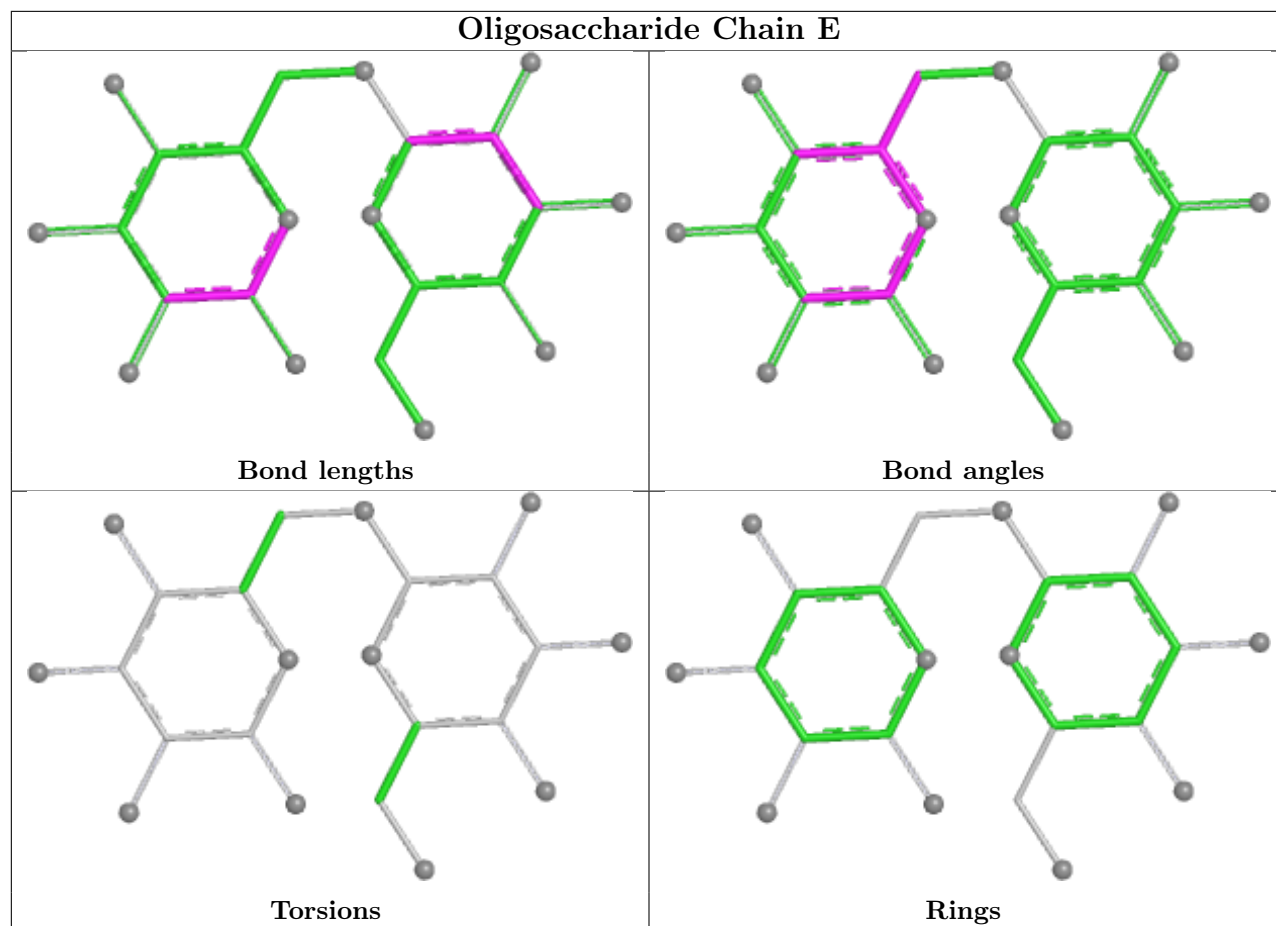
There are no chirality outliers.

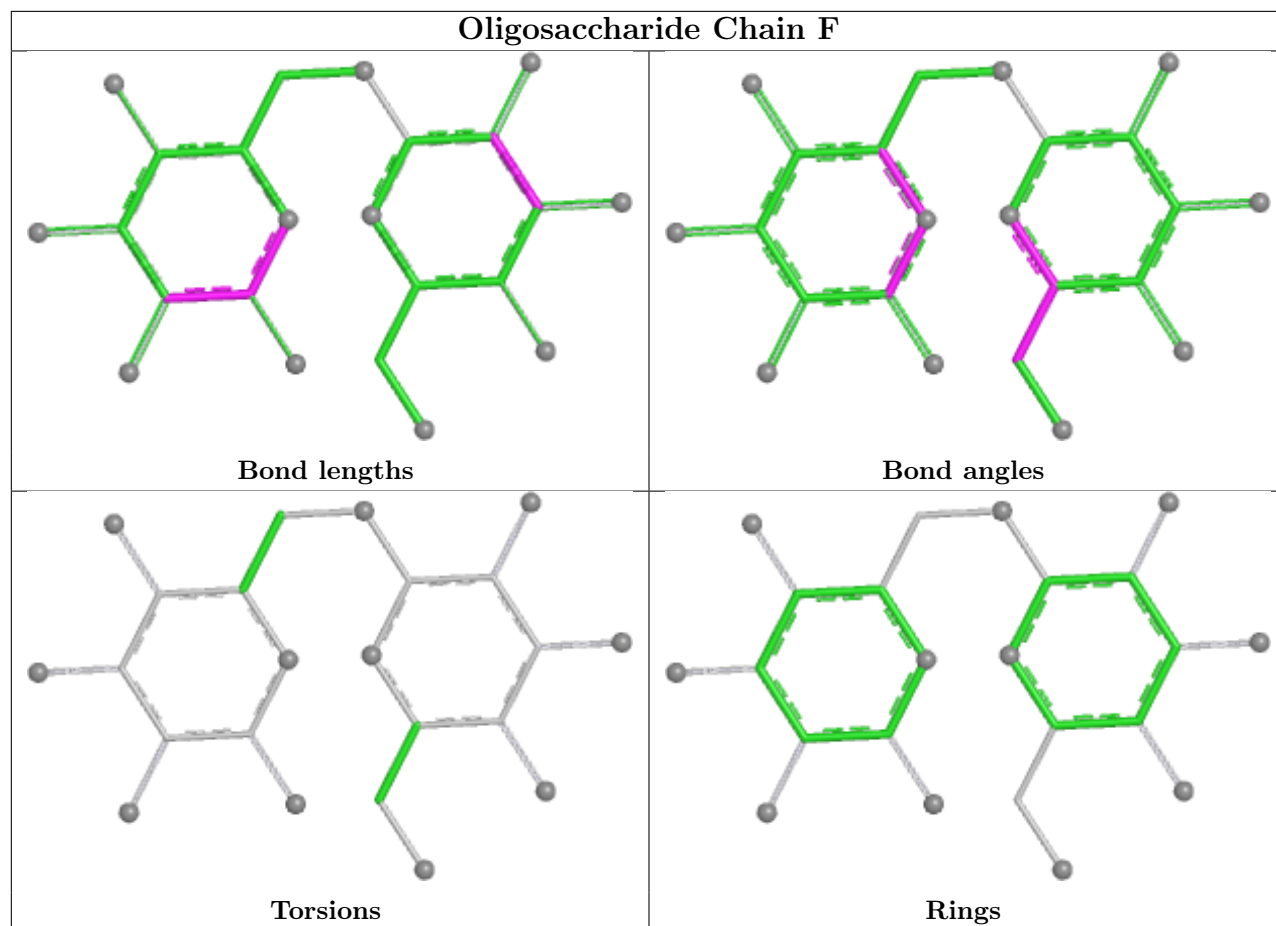
There are no torsion outliers.

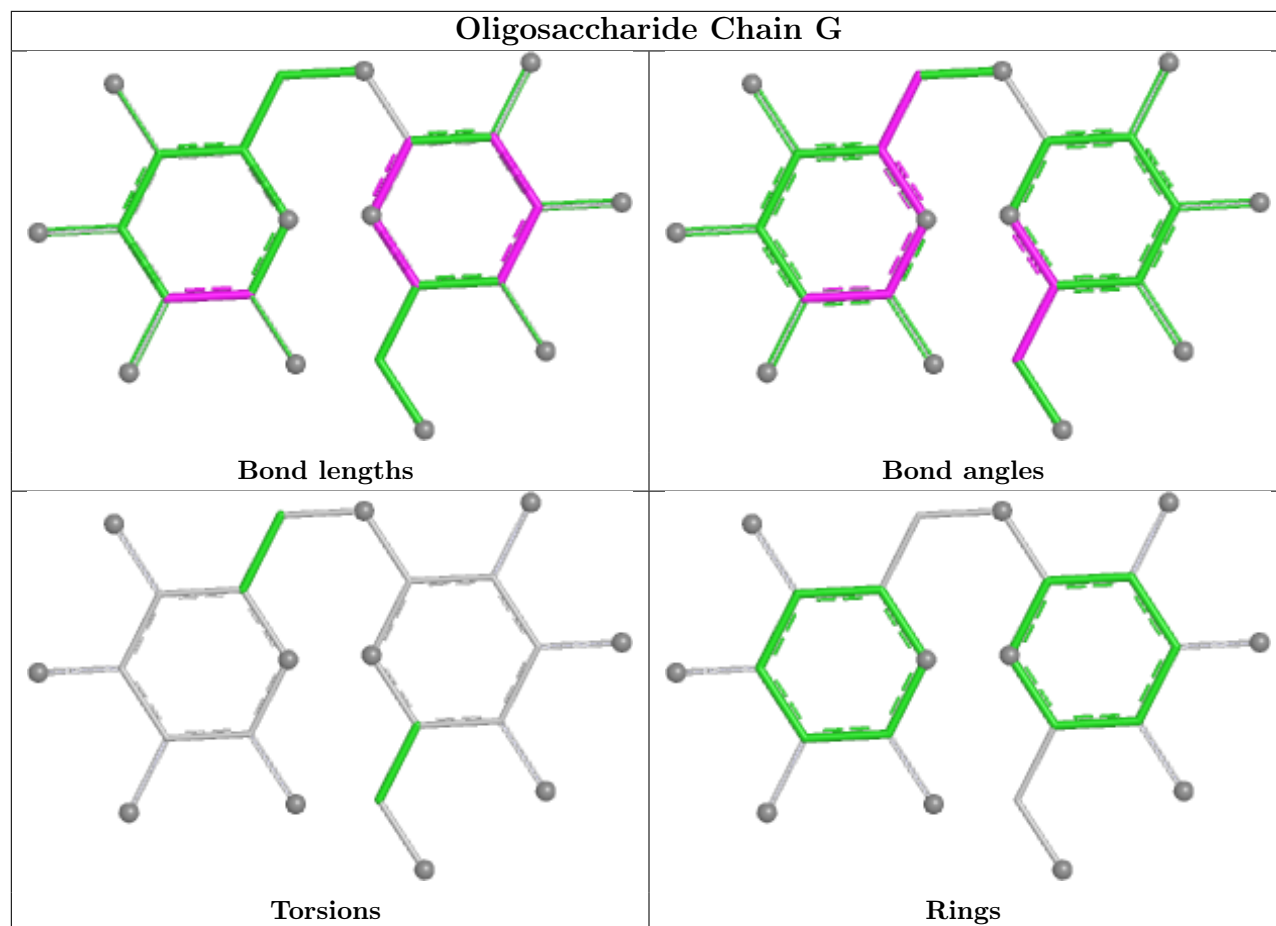
There are no ring outliers.

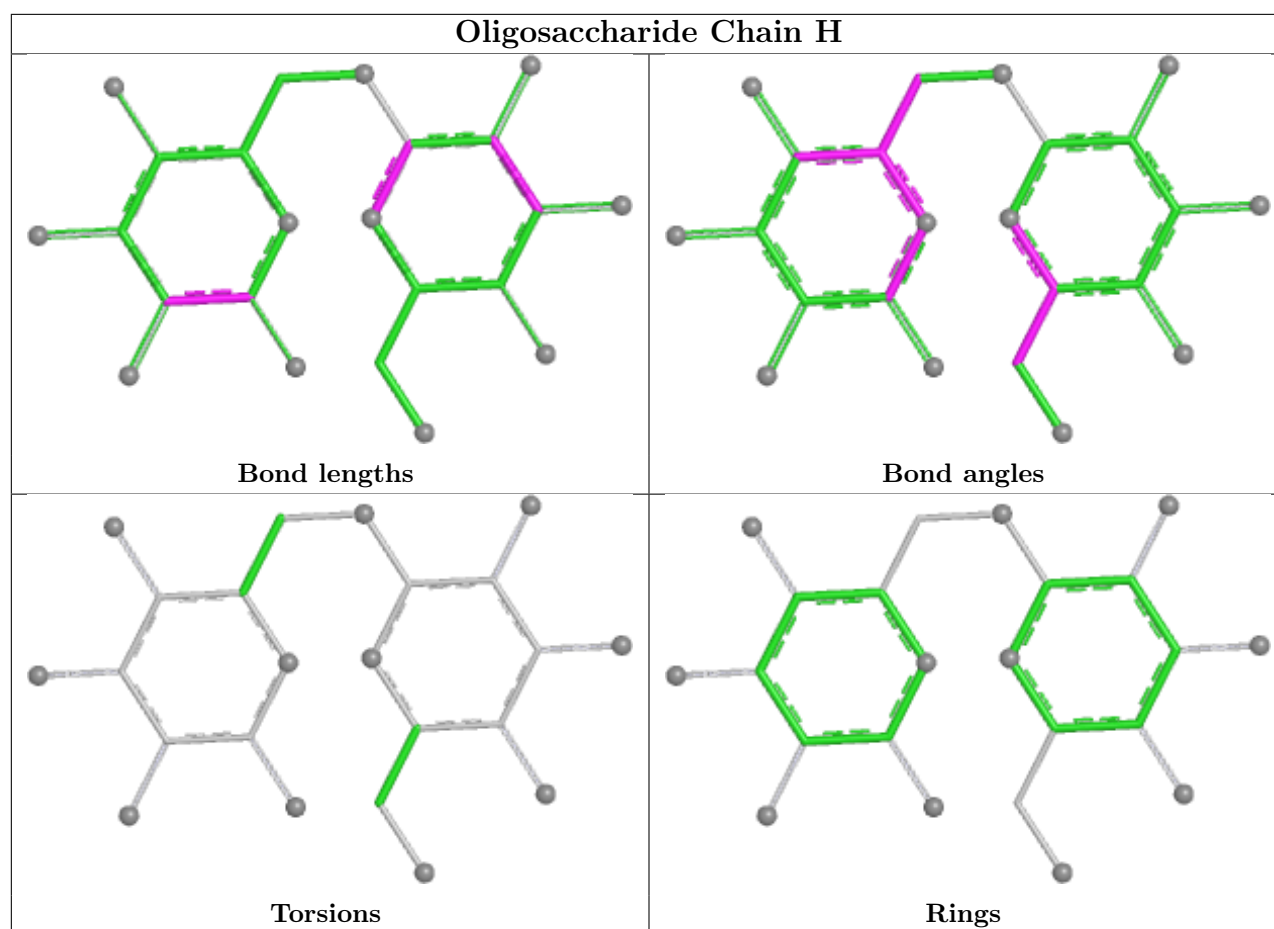
No monomer is involved in short contacts.

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 44 ligands modelled in this entry, 34 are monoatomic - leaving 10 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	$\# Z > 2$
4	EDO	A	705	-	3,3,3	0.55	0	2,2,2	0.44	0
4	EDO	D	703	-	3,3,3	0.53	0	2,2,2	0.31	0
4	EDO	B	703	-	3,3,3	0.45	0	2,2,2	0.50	0
4	EDO	A	703	-	3,3,3	0.61	0	2,2,2	0.36	0
4	EDO	C	703	-	3,3,3	0.59	0	2,2,2	0.21	0
4	EDO	B	702	-	3,3,3	0.68	0	2,2,2	0.16	0
4	EDO	A	704	-	3,3,3	0.47	0	2,2,2	0.39	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
7	PEG	D	710	5	6,6,6	0.24	0	5,5,5	0.33	0
4	EDO	A	706	-	3,3,3	0.63	0	2,2,2	0.07	0
4	EDO	D	702	-	3,3,3	0.54	0	2,2,2	0.25	0

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	EDO	A	705	-	-	1/1/1/1	-
4	EDO	D	703	-	-	0/1/1/1	-
4	EDO	B	703	-	-	1/1/1/1	-
4	EDO	A	703	-	-	0/1/1/1	-
4	EDO	C	703	-	-	0/1/1/1	-
4	EDO	B	702	-	-	0/1/1/1	-
4	EDO	A	704	-	-	0/1/1/1	-
7	PEG	D	710	5	-	2/4/4/4	-
4	EDO	A	706	-	-	1/1/1/1	-
4	EDO	D	702	-	-	1/1/1/1	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (6) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
7	D	710	PEG	O2-C3-C4-O4
4	D	702	EDO	O1-C1-C2-O2
4	A	705	EDO	O1-C1-C2-O2
7	D	710	PEG	C4-C3-O2-C2
4	B	703	EDO	O1-C1-C2-O2
4	A	706	EDO	O1-C1-C2-O2

There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	D	703	EDO	1	0
4	B	702	EDO	1	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

In the following table, the column labelled ‘#RSRZ > 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q < 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	672/683 (98%)	0.00	15 (2%) 62 62	18, 29, 53, 102	0
1	B	671/683 (98%)	0.34	21 (3%) 51 51	19, 35, 64, 100	1 (0%)
1	C	673/683 (98%)	-0.07	9 (1%) 75 75	20, 29, 54, 96	0
1	D	671/683 (98%)	0.08	20 (2%) 52 52	15, 29, 55, 95	1 (0%)
All	All	2687/2732 (98%)	0.09	65 (2%) 59 60	15, 30, 58, 102	2 (0%)

All (65) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	24	GLY	6.8
1	A	22	ALA	5.8
1	D	21	PRO	5.3
1	D	132	ALA	4.8
1	D	24	GLY	4.7
1	D	22	ALA	4.6
1	D	673	GLY	4.0
1	D	23	GLU	3.7
1	A	21	PRO	3.7
1	D	678	HIS	3.5
1	A	564	GLY	3.4
1	C	57	GLY	3.4
1	C	22	ALA	3.2
1	B	670	PRO	3.1
1	A	679	HIS	3.1
1	B	24	GLY	3.0
1	B	671	GLY	3.0
1	D	57	GLY	3.0
1	B	8	ALA	2.9
1	C	24	GLY	2.9
1	B	57	GLY	2.9

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Mol	Chain	Res	Type	RSRZ
1	B	22	ALA	2.8
1	D	133	LYS	2.8
1	D	56	ALA	2.7
1	C	679	HIS	2.7
1	B	463	GLY	2.7
1	B	20	PHE	2.7
1	B	464	PHE	2.6
1	A	351	ASN	2.6
1	B	678	HIS	2.6
1	B	457	LYS	2.6
1	B	504	LEU	2.6
1	C	472	LEU	2.6
1	A	678	HIS	2.5
1	C	574	LYS	2.5
1	A	23	GLU	2.5
1	B	467	PHE	2.5
1	B	21	PRO	2.5
1	D	55	LYS	2.5
1	C	678	HIS	2.5
1	B	472	LEU	2.4
1	D	674	GLY	2.4
1	D	20	PHE	2.4
1	C	587	GLN	2.4
1	B	551	VAL	2.4
1	A	20	PHE	2.3
1	B	511	HIS	2.3
1	A	350	PRO	2.3
1	D	236	CYS	2.3
1	D	52	LYS	2.2
1	C	20	PHE	2.2
1	A	133	LYS	2.2
1	D	471	THR	2.2
1	D	350	PRO	2.2
1	A	565	LYS	2.2
1	B	56	ALA	2.2
1	B	503	LYS	2.2
1	B	554	ARG	2.2
1	A	514	ASP	2.1
1	B	58	LYS	2.1
1	D	472	LEU	2.0
1	D	554	ARG	2.0
1	A	563	GLU	2.0

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Mol	Chain	Res	Type	RSRZ
1	D	511	HIS	2.0
1	A	562	PRO	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](#)

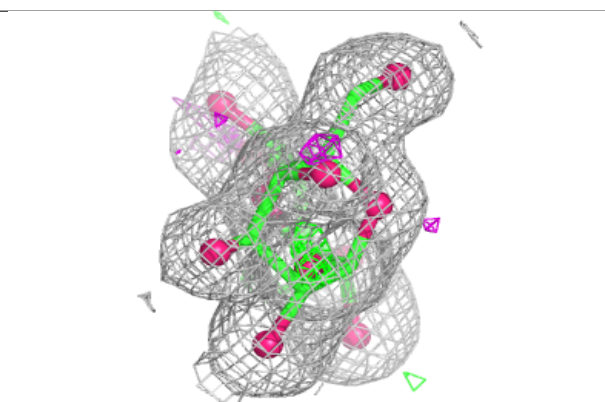
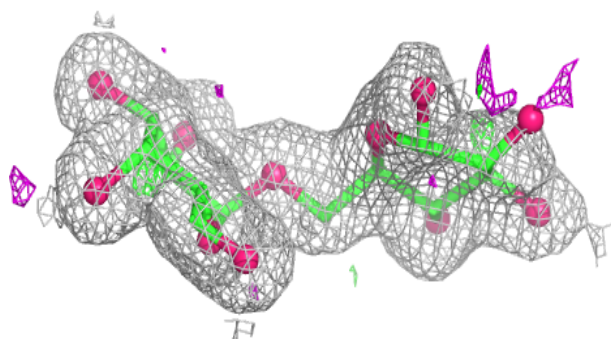
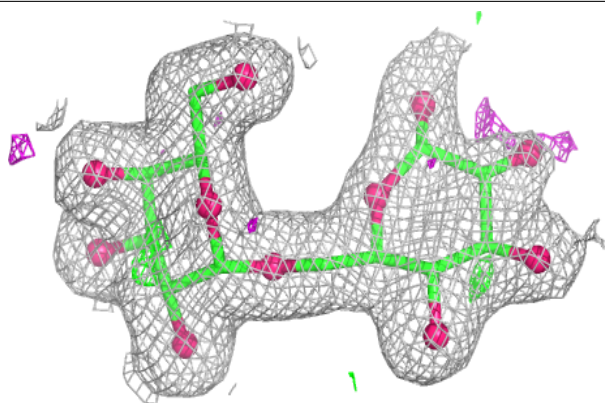
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q<0.9' lists 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	E	1	12/12	0.89	0.12	23,40,49,53	0
2	GLC	H	1	12/12	0.90	0.11	25,40,47,48	0
2	GLC	G	1	12/12	0.93	0.09	22,35,44,50	0
2	GLC	F	1	12/12	0.93	0.09	25,37,49,50	0
2	GLA	E	2	11/12	0.96	0.06	18,20,21,22	0
2	GLA	G	2	11/12	0.97	0.06	18,19,22,23	0
2	GLA	F	2	11/12	0.97	0.05	20,21,22,24	0
2	GLA	H	2	11/12	0.97	0.05	18,23,27,27	0

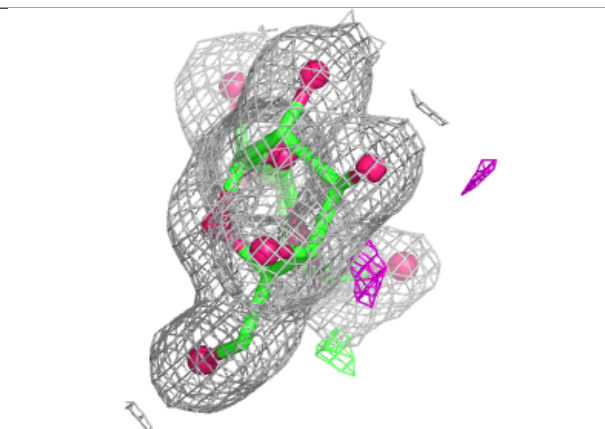
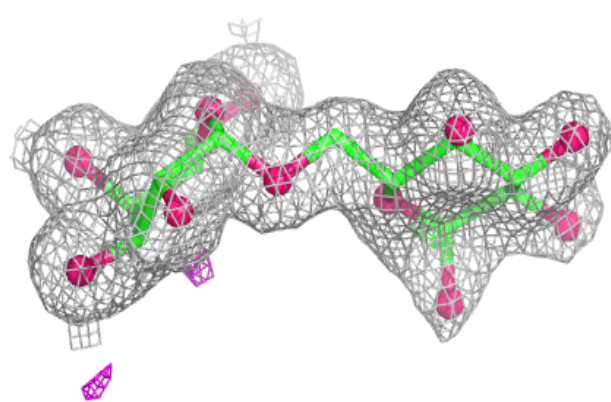
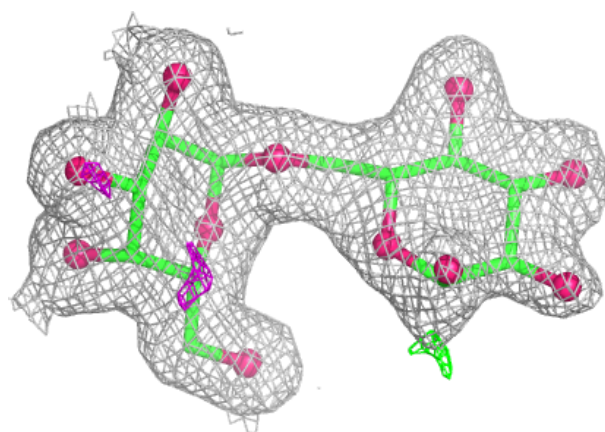
The following is a graphical depiction of the model fit to experimental electron density for oligosaccharide. Each fit is shown from different orientation to approximate a three-dimensional view.

Electron density around Chain E:

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

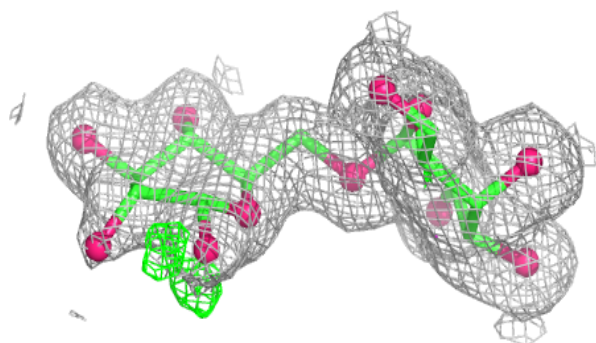
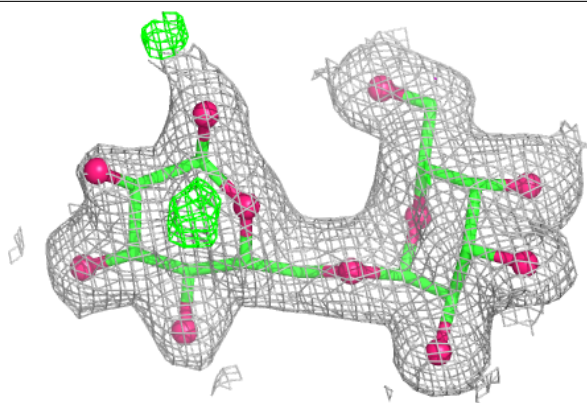
**Electron density around Chain F:**

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

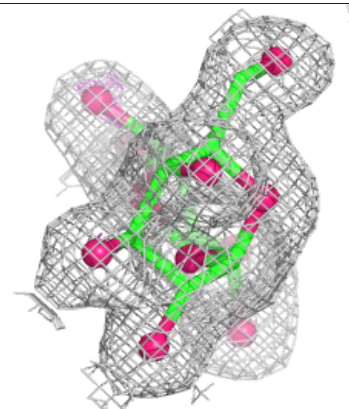
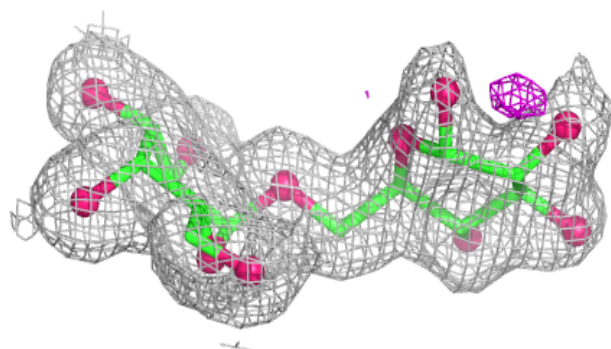
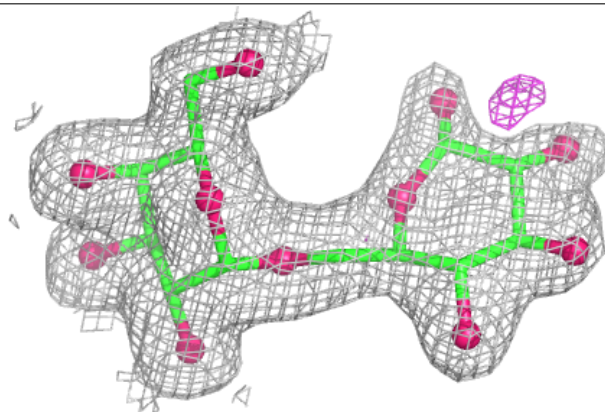


Electron density around Chain G:

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

**Electron density around Chain H:**

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



6.4 Ligands ⓘ

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	EDO	A	705	4/4	0.48	0.30	62,64,64,65	0
4	EDO	A	706	4/4	0.75	0.18	70,70,70,70	0
4	EDO	B	702	4/4	0.77	0.16	42,44,45,47	0
7	PEG	D	710	7/7	0.79	0.14	40,42,47,48	0
4	EDO	D	703	4/4	0.81	0.15	65,65,65,66	0
4	EDO	A	704	4/4	0.82	0.14	61,61,62,63	0
4	EDO	C	703	4/4	0.82	0.14	38,41,43,44	0
4	EDO	B	703	4/4	0.83	0.14	38,47,51,56	0
5	NA	A	708	1/1	0.84	0.17	55,55,55,55	0
5	NA	A	712	1/1	0.88	0.20	63,63,63,63	0
5	NA	C	712	1/1	0.90	0.12	40,40,40,40	0
5	NA	B	713	1/1	0.91	0.13	39,39,39,39	0
4	EDO	A	703	4/4	0.91	0.08	32,34,36,36	0
5	NA	C	710	1/1	0.92	0.11	47,47,47,47	0
4	EDO	D	702	4/4	0.92	0.10	57,58,59,60	0
5	NA	A	709	1/1	0.93	0.16	51,51,51,51	0
5	NA	D	705	1/1	0.94	0.12	35,35,35,35	0
5	NA	B	707	1/1	0.94	0.07	36,36,36,36	0
5	NA	B	709	1/1	0.95	0.07	38,38,38,38	0
5	NA	B	712	1/1	0.95	0.09	42,42,42,42	0
5	NA	D	707	1/1	0.96	0.07	49,49,49,49	0
5	NA	C	709	1/1	0.96	0.07	42,42,42,42	0
3	MG	B	701	1/1	0.96	0.07	39,39,39,39	0
3	MG	D	701	1/1	0.96	0.06	41,41,41,41	0
5	NA	C	713	1/1	0.96	0.06	41,41,41,41	0
3	MG	C	701	1/1	0.97	0.05	29,29,29,29	0
5	NA	D	708	1/1	0.97	0.05	35,35,35,35	0
5	NA	C	706	1/1	0.97	0.07	37,37,37,37	0
5	NA	A	710	1/1	0.97	0.07	34,34,34,34	0
5	NA	A	707	1/1	0.98	0.04	27,27,27,27	0
5	NA	A	711	1/1	0.98	0.05	33,33,33,33	0
5	NA	B	705	1/1	0.98	0.07	24,24,24,24	0
5	NA	B	706	1/1	0.98	0.07	32,32,32,32	0
5	NA	C	705	1/1	0.98	0.06	25,25,25,25	0
5	NA	B	708	1/1	0.98	0.10	26,26,26,26	0
3	MG	A	702	1/1	0.98	0.07	28,28,28,28	0
5	NA	B	710	1/1	0.98	0.04	35,35,35,35	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
5	NA	B	711	1/1	0.98	0.10	37,37,37,37	0
3	MG	C	702	1/1	0.98	0.04	36,36,36,36	0
5	NA	D	706	1/1	0.98	0.06	29,29,29,29	0
6	CL	D	709	1/1	0.98	0.11	37,37,37,37	0
5	NA	C	711	1/1	0.98	0.04	27,27,27,27	0
5	NA	C	707	1/1	0.99	0.05	31,31,31,31	0
5	NA	C	708	1/1	0.99	0.08	24,24,24,24	0

6.5 Other polymers [i](#)

There are no such residues in this entry.