



Full wwPDB X-ray Structure Validation Report ⓘ

May 28, 2025 – 10:12 PM JST

PDB ID : 9KYQ / pdb_00009kyq
Title : GH57 family amylopullulanase from Aquifex aeolicus wild type co-crystallize with gama-cyclodextrin
Authors : Zhu, Z.M.; Wang, W.W.; Yu, F.; Li, M.J.; Xu, Q.; Zhou, H.; Huang, L.Q.; Wang, Q.S.
Deposited on : 2024-12-09
Resolution : 1.60 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 Phenix2.0rc1
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 2.0rc1
EDS : 3.0
Percentile statistics : 20231227.v01 (using entries in the PDB archive December 27th 2023)
CCP4 : 9.0.006 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.43.1

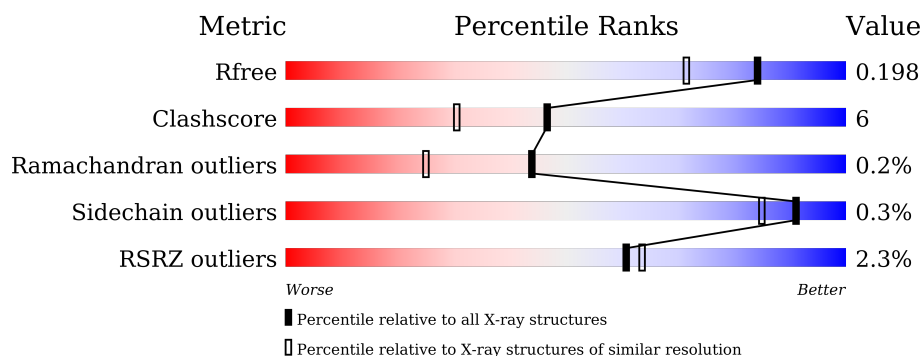
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.60 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}	164625	4274 (1.60-1.60)
Clashscore	180529	4682 (1.60-1.60)
Ramachandran outliers	177936	4583 (1.60-1.60)
Sidechain outliers	177891	4582 (1.60-1.60)
RSRZ outliers	164620	4272 (1.60-1.60)

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	477	<div> <div style="width: 100%; height: 10px; background-color: red;"></div> <div style="width: 90%; height: 10px; background-color: green;"></div> <div style="width: 10%; height: 10px; background-color: yellow;"></div> <div>90%10%</div> </div>
2	B	478	<div> <div style="width: 4%; height: 10px; background-color: red;"></div> <div style="width: 88%; height: 10px; background-color: green;"></div> <div style="width: 12%; height: 10px; background-color: yellow;"></div> <div>4%88%12%</div> </div>
3	C	3	<div> <div style="width: 33%; height: 10px; background-color: green;"></div> <div style="width: 33%; height: 10px; background-color: yellow;"></div> <div style="width: 33%; height: 10px; background-color: orange;"></div> <div>33%33%33%</div> </div>
3	D	3	<div> <div style="width: 100%; height: 10px; background-color: yellow;"></div> <div>100%</div> </div>

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit crite-

ria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
6	PEG	A	506	-	-	X	-

2 Entry composition [i](#)

There are 7 unique types of molecules in this entry. The entry contains 8956 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 Glycoside hydrolase family 57 N-terminal domain-containing protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	476	Total	C	N	O	S	0	5	0
			4065	2659	652	744	10			

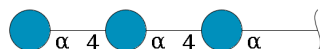
- Molecule 2 is a protein called Glycoside hydrolase family 57 N-terminal domain-containing protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	478	Total	C	N	O	S	0	3	0
			4058	2656	649	742	11			

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	478	HIS	-	expression tag	UNP O66934

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



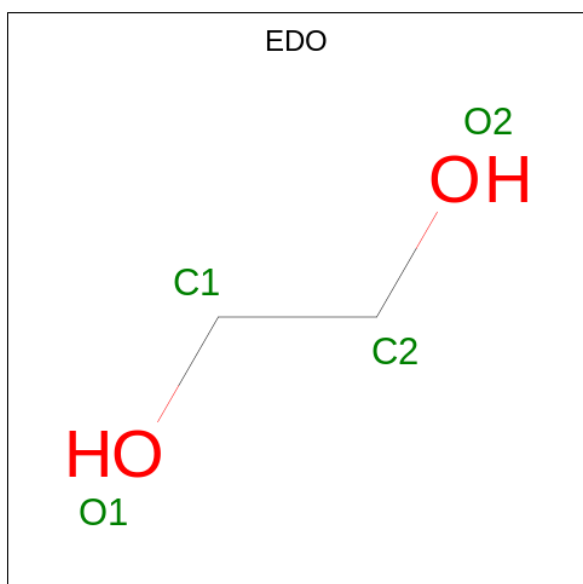
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace
3	C	3	Total	C	O	0	0	0
			34	18	16			
3	D	3	Total	C	O	0	0	0
			34	18	16			

- Molecule 4 is GLYCEROL (CCD ID: GOL) (formula: C₃H₈O₃).



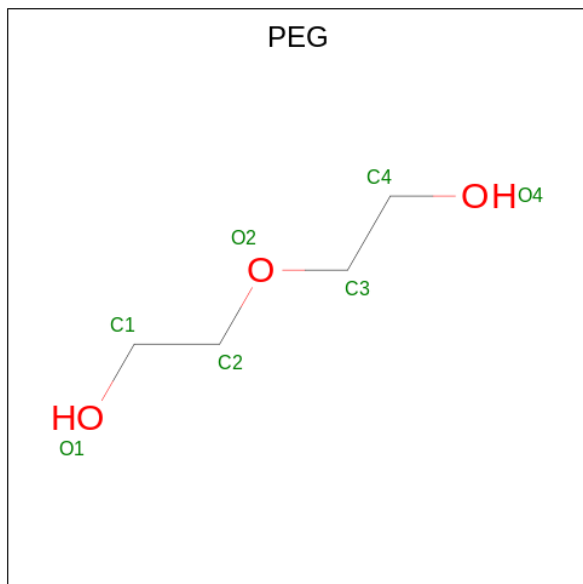
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			6	3	3		
4	A	1	Total	C	O	0	0
			6	3	3		
4	A	1	Total	C	O	0	0
			6	3	3		
4	B	1	Total	C	O	0	0
			6	3	3		
4	B	1	Total	C	O	0	0
			6	3	3		

- Molecule 5 is 1,2-ETHANEDIOL (CCD ID: EDO) (formula: C₂H₆O₂).



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

- Molecule 6 is DI(HYDROXYETHYL)ETHER (CCD ID: PEG) (formula: C₄H₁₀O₃).



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

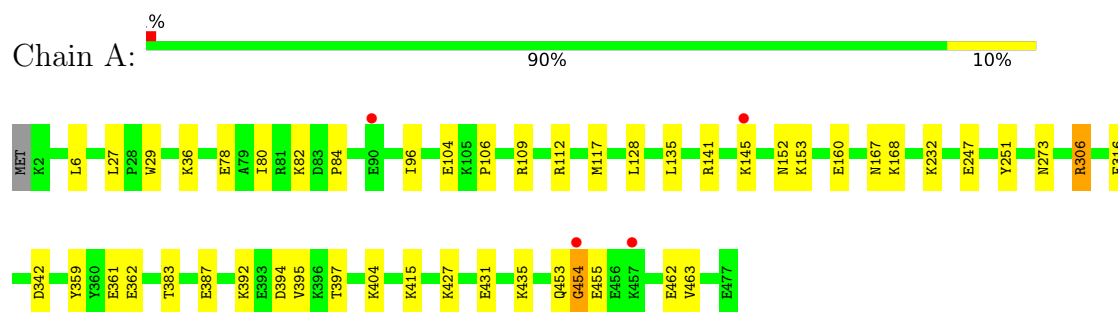
- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	406	Total	O	0	0
			406	406		
7	B	310	Total	O	0	0
			310	310		

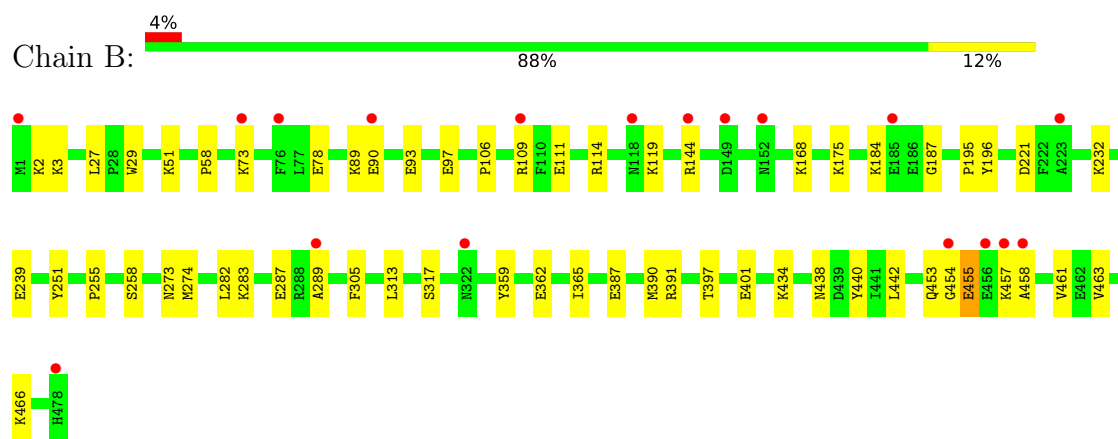
3 Residue-property plots [i](#)

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.

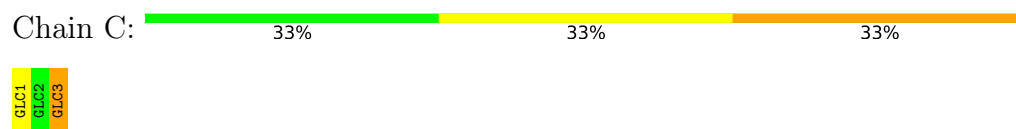
- Molecule 1: Glycoside hydrolase family 57 N-terminal domain-containing protein



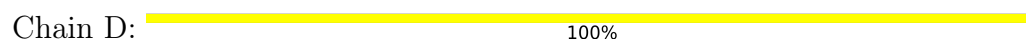
- Molecule 2: Glycoside hydrolase family 57 N-terminal domain-containing protein



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



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



GLC1
GLC2
GLC3

4 Data and refinement statistics

Property	Value	Source
Space group	P 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	61.45Å 41.35Å 194.02Å 90.00° 96.24° 90.00°	Depositor
Resolution (Å)	64.29 – 1.60 64.29 – 1.60	Depositor EDS
% Data completeness (in resolution range)	99.5 (64.29-1.60) 99.5 (64.29-1.60)	Depositor EDS
R_{merge}	0.13	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.72 (at 1.60Å)	Xtriage
Refinement program	PHENIX (1.13_2998: ???)	Depositor
R, R_{free}	0.170 , 0.199 0.170 , 0.198	Depositor DCC
R_{free} test set	6285 reflections (4.87%)	wwPDB-VP
Wilson B-factor (Å ²)	18.2	Xtriage
Anisotropy	0.486	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 42.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.015 for h,-k,-h-l	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	8956	wwPDB-VP
Average B, all atoms (Å ²)	25.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 11.35% 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: GOL, GLC, EDO, PEG

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	0.36	1/4179 (0.0%)	0.54	2/5656 (0.0%)
2	B	0.35	3/4172 (0.1%)	0.55	1/5647 (0.0%)
All	All	0.36	4/8351 (0.0%)	0.54	3/11303 (0.0%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1
2	B	0	1
All	All	0	2

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	2	LYS	C-O	-7.34	1.15	1.24
1	A	306	ARG	C-O	-6.69	1.15	1.23
2	B	2	LYS	CA-CB	-5.57	1.45	1.53
2	B	3	LYS	C-O	-5.47	1.17	1.23

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	2	LYS	N-CA-C	-14.30	85.21	108.52
1	A	454	GLY	CA-C-N	5.40	131.41	121.70
1	A	454	GLY	C-N-CA	5.40	131.41	121.70

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	453	GLN	Peptide
2	B	453	GLN	Peptide

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	4065	0	3967	39	0
2	B	4058	0	3964	48	0
3	C	34	0	29	1	0
3	D	34	0	30	1	0
4	A	18	0	23	3	0
4	B	12	0	15	2	0
5	A	8	0	12	0	0
5	B	4	0	6	0	0
6	A	7	0	10	7	0
7	A	406	0	0	16	2
7	B	310	0	0	15	2
All	All	8956	0	8056	91	3

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 6.

All (91) 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)
2:B:144:ARG:NH1	7:B:601:HOH:O	1.85	1.01
1:A:316:PHE:CE2	6:A:506:PEG:H12	1.97	0.99
1:A:316:PHE:HE2	6:A:506:PEG:H12	1.33	0.92
2:B:251:TYR:HE1	2:B:390:MET:HE3	1.34	0.91
2:B:144:ARG:NH2	7:B:603:HOH:O	2.03	0.90
2:B:434:LYS:NZ	7:B:602:HOH:O	1.98	0.81
1:A:152:ASN:OD1	7:A:601:HOH:O	2.05	0.74
2:B:362:GLU:OE1	7:B:604:HOH:O	2.05	0.72
2:B:287:GLU:OE2	7:B:605:HOH:O	2.10	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:251:TYR:CE1	2:B:390:MET:HE3	2.23	0.70
2:B:119:LYS:NZ	7:B:611:HOH:O	2.24	0.69
1:A:361:GLU:HG2	1:A:362:GLU:HG3	1.74	0.69
2:B:221:ASP:OD2	7:B:607:HOH:O	2.13	0.66
1:A:160[B]:GLU:OE1	7:A:602:HOH:O	2.13	0.65
2:B:401:GLU:OE2	7:B:606:HOH:O	2.13	0.65
6:A:506:PEG:H11	7:A:882:HOH:O	1.98	0.64
2:B:29:TRP:CD2	3:D:3:GLC:H2	2.37	0.60
1:A:106:PRO:HA	1:A:109:ARG:HG3	1.85	0.59
1:A:29:TRP:CD2	3:C:3:GLC:H2	2.38	0.59
2:B:273:ASN:HB2	2:B:390:MET:HE1	1.85	0.59
1:A:247:GLU:OE1	7:A:603:HOH:O	2.16	0.59
1:A:454:GLY:HA2	7:A:662:HOH:O	2.04	0.57
1:A:104:GLU:OE2	7:A:604:HOH:O	2.18	0.57
2:B:73:LYS:NZ	2:B:78:GLU:HG3	2.20	0.55
2:B:90:GLU:HA	2:B:93:GLU:HG3	1.88	0.55
2:B:109:ARG:HG2	7:B:672:HOH:O	2.06	0.55
2:B:282:LEU:HG	2:B:289:ALA:HB1	1.87	0.55
1:A:431:GLU:CD	1:A:435:LYS:HZ3	2.14	0.54
2:B:184:LYS:HZ3	4:B:502:GOL:H31	1.71	0.54
1:A:167:ASN:HB3	7:A:841:HOH:O	2.08	0.53
2:B:239:GLU:OE2	7:B:610:HOH:O	2.19	0.53
1:A:112:ARG:HH22	4:A:503:GOL:H2	1.74	0.53
2:B:391:ARG:NH2	7:B:612:HOH:O	2.27	0.52
2:B:387:GLU:OE2	2:B:391:ARG:NH2	2.42	0.52
2:B:89:LYS:O	2:B:93:GLU:HG3	2.09	0.52
1:A:394:ASP:OD1	1:A:394:ASP:N	2.38	0.51
2:B:387:GLU:HG3	2:B:391:ARG:HE	1.75	0.51
1:A:404:LYS:HE3	7:A:814:HOH:O	2.11	0.51
2:B:195:PRO:HD2	2:B:255:PRO:HD3	1.93	0.50
2:B:258[A]:SER:OG	7:B:609:HOH:O	2.17	0.50
2:B:184:LYS:NZ	4:B:502:GOL:H31	2.28	0.48
1:A:306:ARG:HH12	6:A:506:PEG:H21	1.77	0.48
2:B:274[B]:MET:HB2	2:B:390:MET:HE2	1.95	0.48
2:B:282:LEU:HB2	2:B:305:PHE:CD2	2.49	0.47
1:A:112:ARG:NH2	4:A:503:GOL:H2	2.29	0.47
2:B:283:LYS:HD3	2:B:289:ALA:HB3	1.97	0.47
1:A:273:ASN:ND2	7:A:620:HOH:O	2.47	0.47
1:A:84:PRO:HG2	1:A:128:LEU:HG	1.97	0.47
2:B:438:ASN:OD1	2:B:440:TYR:HB2	2.15	0.47
1:A:247:GLU:HB3	7:A:603:HOH:O	2.16	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:232:LYS:HB3	1:A:232:LYS:HE3	1.45	0.46
2:B:106:PRO:HA	2:B:109:ARG:HG3	1.98	0.46
2:B:144:ARG:CZ	2:B:365:ILE:HD13	2.46	0.45
6:A:506:PEG:C1	7:A:882:HOH:O	2.60	0.45
2:B:73:LYS:HD2	2:B:73:LYS:HA	1.72	0.45
2:B:73:LYS:HZ2	2:B:78:GLU:HG3	1.80	0.45
1:A:106:PRO:HD2	1:A:359:TYR:HA	1.98	0.44
1:A:80:ILE:HG21	1:A:135:LEU:HG	1.99	0.44
2:B:458:ALA:HB1	2:B:461:VAL:HB	2.00	0.44
1:A:462:GLU:HG3	7:A:875:HOH:O	2.17	0.44
1:A:168:LYS:HE3	7:A:849:HOH:O	2.17	0.44
1:A:27:LEU:HG	1:A:463:VAL:HG11	1.99	0.44
2:B:251:TYR:OH	2:B:387:GLU:HA	2.18	0.44
2:B:313:LEU:HA	2:B:317:SER:HB2	2.00	0.43
1:A:342:ASP:OD1	1:A:342:ASP:N	2.42	0.43
2:B:93:GLU:O	2:B:97:GLU:HG3	2.18	0.43
2:B:106:PRO:HD2	2:B:359:TYR:HA	1.99	0.43
2:B:168:LYS:HE3	7:B:775:HOH:O	2.18	0.43
1:A:427:LYS:HG3	7:A:812:HOH:O	2.18	0.42
1:A:153:LYS:HE3	1:A:153:LYS:HB3	1.77	0.42
2:B:232:LYS:HD2	2:B:442:LEU:HD11	2.01	0.42
1:A:454:GLY:HA2	1:A:455:GLU:HB2	2.01	0.42
1:A:392:LYS:HA	1:A:392:LYS:HD3	1.90	0.41
2:B:454:GLY:HA2	2:B:455:GLU:HB2	2.03	0.41
2:B:387:GLU:CG	2:B:391:ARG:HE	2.33	0.41
6:A:506:PEG:H22	7:A:882:HOH:O	2.20	0.41
2:B:466:LYS:NZ	7:B:623:HOH:O	2.51	0.41
2:B:58:PRO:HG2	2:B:196:TYR:HA	2.02	0.41
2:B:175:LYS:HG3	7:B:786:HOH:O	2.20	0.41
1:A:78:GLU:O	1:A:82:LYS:HG3	2.20	0.41
1:A:251:TYR:OH	1:A:387:GLU:HA	2.21	0.41
6:A:506:PEG:C1	6:A:506:PEG:H42	2.49	0.41
2:B:111:GLU:OE2	2:B:114:ARG:NH2	2.54	0.40
2:B:51:LYS:NZ	2:B:187:GLY:O	2.51	0.40
1:A:167:ASN:CB	7:A:841:HOH:O	2.67	0.40
1:A:415:LYS:HD2	1:A:415:LYS:HA	1.86	0.40
2:B:27:LEU:HG	2:B:463:VAL:HG11	2.04	0.40
1:A:96:ILE:HG23	1:A:117:MET:HE1	2.03	0.40
1:A:6:LEU:O	1:A:383:THR:HA	2.21	0.40
1:A:36:LYS:O	4:A:502:GOL:H32	2.21	0.40
1:A:141:ARG:O	1:A:145:LYS:HD2	2.21	0.40

All (3) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
7:A:654:HOH:O	7:A:850:HOH:O[1_455]	1.97	0.23
7:A:790:HOH:O	7:B:653:HOH:O[1_645]	2.07	0.13
7:B:668:HOH:O	7:B:692:HOH:O[1_455]	2.09	0.11

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	479/477 (100%)	472 (98%)	7 (2%)	0	100	100
2	B	479/478 (100%)	469 (98%)	8 (2%)	2 (0%)	30	14
All	All	958/955 (100%)	941 (98%)	15 (2%)	2 (0%)	44	25

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	457	LYS
2	B	455	GLU

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	435/437 (100%)	433 (100%)	2 (0%)	86	78
2	B	434/438 (99%)	433 (100%)	1 (0%)	92	86

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
All	All	869/875 (99%)	866 (100%)	3 (0%)	91	85

All (3) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	395	VAL
1	A	397	THR
2	B	397	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	118	ASN
1	A	156	ASN
1	A	476	GLN
2	B	476	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 ⓘ

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	$\# Z > 2$
3	GLC	C	1	3	12,12,12	1.06	1 (8%)	17,17,17	1.44	3 (17%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	GLC	C	2	3	11,11,12	0.47	0	15,15,17	0.90	0
3	GLC	C	3	3	11,11,12	0.48	0	15,15,17	1.06	1 (6%)
3	GLC	D	1	3	12,12,12	1.52	2 (16%)	17,17,17	1.52	2 (11%)
3	GLC	D	2	3	11,11,12	0.66	0	15,15,17	1.00	1 (6%)
3	GLC	D	3	3	11,11,12	0.47	0	15,15,17	0.93	0

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	GLC	C	1	3	-	0/2/22/22	0/1/1/1
3	GLC	C	2	3	-	0/2/19/22	0/1/1/1
3	GLC	C	3	3	-	0/2/19/22	0/1/1/1
3	GLC	D	1	3	-	0/2/22/22	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

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	D	1	GLC	O5-C5	-3.53	1.35	1.44
3	C	1	GLC	O5-C5	-2.95	1.37	1.44
3	D	1	GLC	O5-C1	-2.62	1.36	1.42

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	D	1	GLC	O5-C5-C4	3.84	116.67	109.69
3	D	1	GLC	C1-O5-C5	3.54	120.34	113.66
3	C	1	GLC	O3-C3-C2	-3.36	102.57	110.35
3	C	1	GLC	C1-O5-C5	2.67	118.69	113.66
3	C	1	GLC	O5-C5-C4	2.49	114.22	109.69
3	C	3	GLC	C1-O5-C5	2.46	115.52	112.19
3	D	2	GLC	C1-O5-C5	2.08	115.00	112.19

There are no chirality outliers.

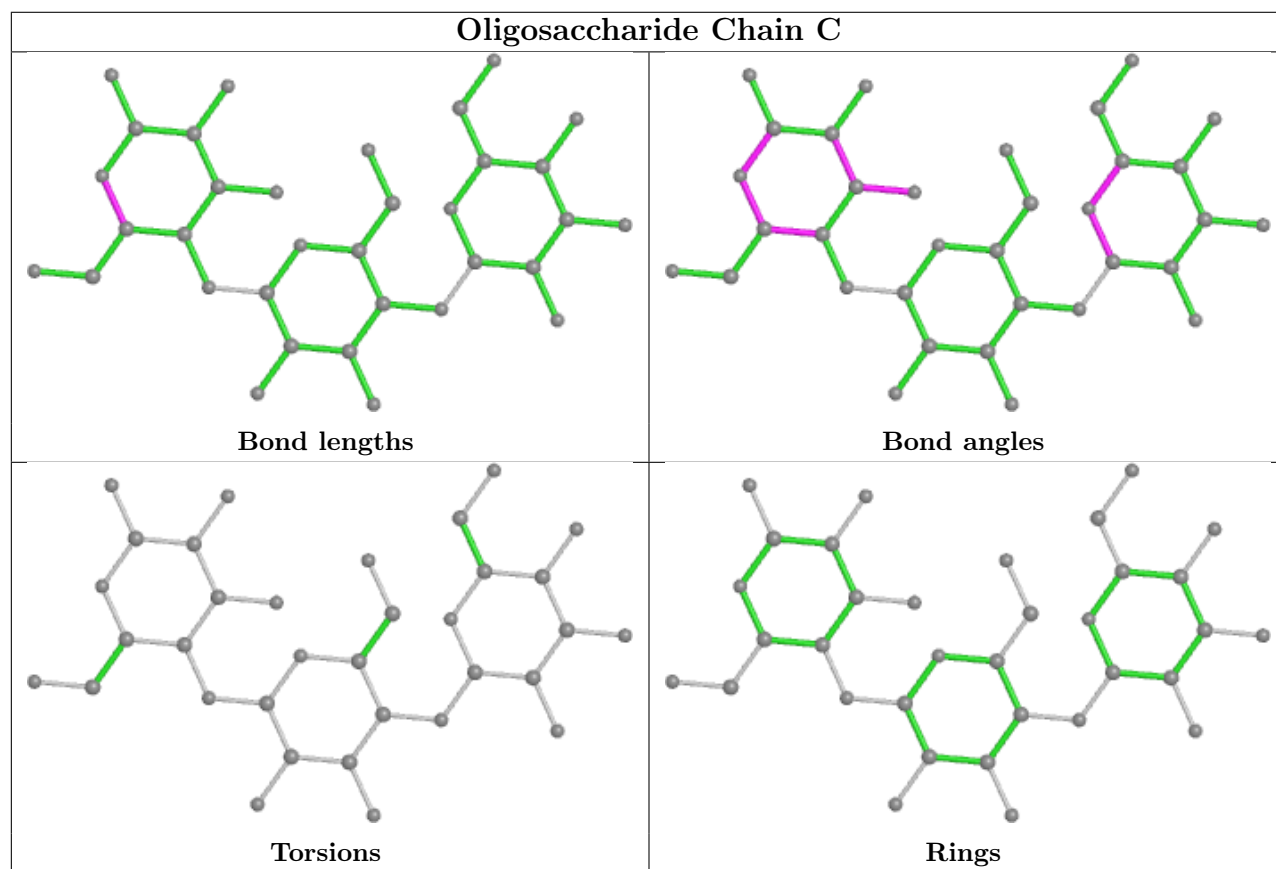
There are no torsion outliers.

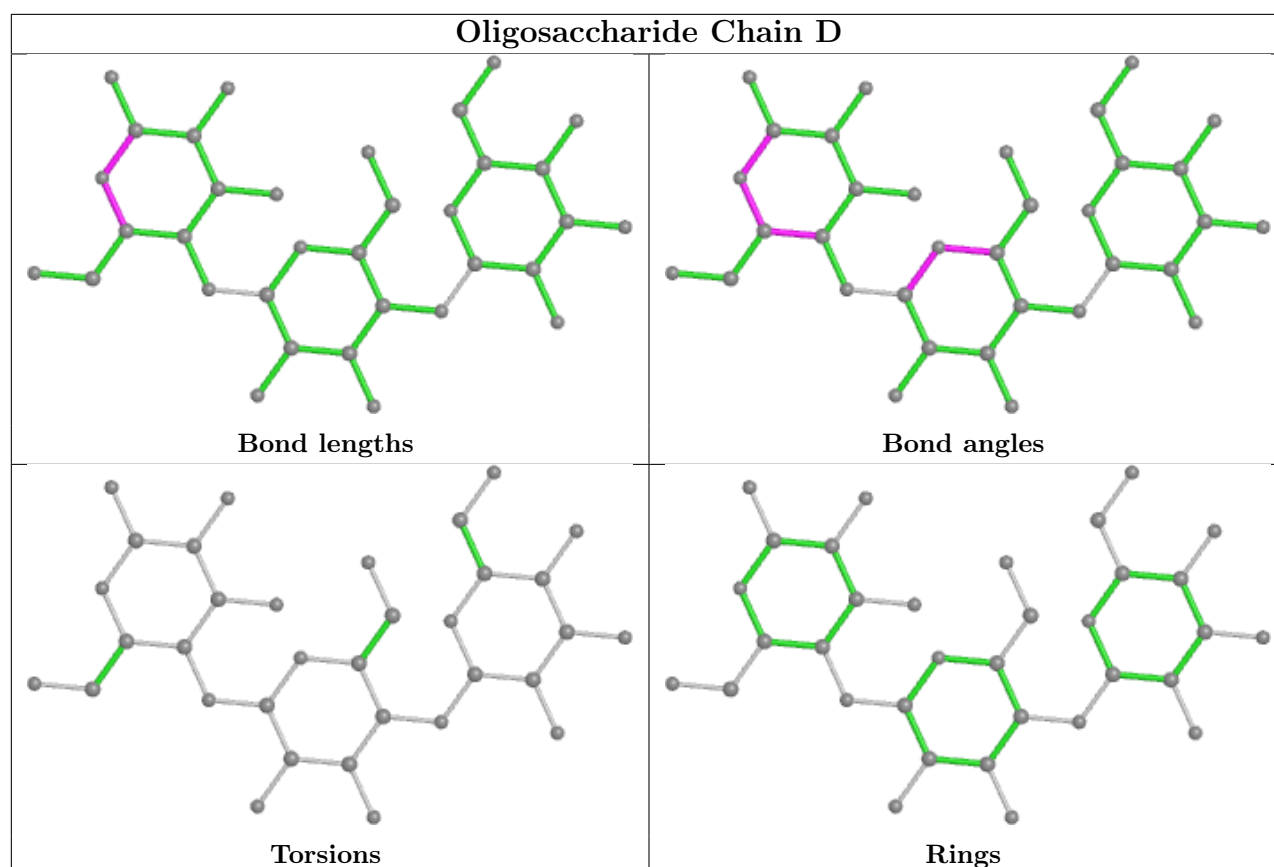
There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	D	3	GLC	1	0
3	C	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](#)

9 ligands are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
4	GOL	B	501	-	5,5,5	1.03	0	5,5,5	0.97	0
4	GOL	A	502	-	5,5,5	1.70	1 (20%)	5,5,5	1.66	1 (20%)
6	PEG	A	506	-	6,6,6	0.40	0	5,5,5	1.19	0
5	EDO	B	503	-	3,3,3	0.43	0	2,2,2	0.51	0
5	EDO	A	504	-	3,3,3	0.45	0	2,2,2	0.35	0
4	GOL	A	503	-	5,5,5	0.75	0	5,5,5	1.39	1 (20%)
5	EDO	A	505	-	3,3,3	0.52	0	2,2,2	0.33	0
4	GOL	A	501	-	5,5,5	1.01	0	5,5,5	0.91	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
4	GOL	B	502	-	5,5,5	1.09	0	5,5,5	3.09	1 (20%)

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	GOL	B	501	-	-	0/4/4/4	-
4	GOL	A	502	-	-	0/4/4/4	-
6	PEG	A	506	-	-	3/4/4/4	-
5	EDO	B	503	-	-	1/1/1/1	-
5	EDO	A	504	-	-	0/1/1/1	-
4	GOL	A	503	-	-	2/4/4/4	-
5	EDO	A	505	-	-	0/1/1/1	-
4	GOL	A	501	-	-	0/4/4/4	-
4	GOL	B	502	-	-	2/4/4/4	-

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	502	GOL	O3-C3	-3.42	1.28	1.42

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	502	GOL	O3-C3-C2	-6.84	77.40	110.20
4	A	502	GOL	C3-C2-C1	-2.91	100.39	111.70
4	A	503	GOL	O1-C1-C2	-2.24	99.46	110.20

There are no chirality outliers.

All (8) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	B	502	GOL	O2-C2-C3-O3
6	A	506	PEG	C4-C3-O2-C2
6	A	506	PEG	O1-C1-C2-O2
4	B	502	GOL	C1-C2-C3-O3
6	A	506	PEG	O2-C3-C4-O4
4	A	503	GOL	O1-C1-C2-O2

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Mol	Chain	Res	Type	Atoms
5	B	503	EDO	O1-C1-C2-O2
4	A	503	GOL	O1-C1-C2-C3

There are no ring outliers.

4 monomers are involved in 12 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	502	GOL	1	0
6	A	506	PEG	7	0
4	A	503	GOL	2	0
4	B	502	GOL	2	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 [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	476/477 (99%)	-0.20	4 (0%) 82 85	7, 19, 37, 52	5 (1%)
2	B	478/478 (100%)	0.10	18 (3%) 44 44	8, 25, 45, 71	3 (0%)
All	All	954/955 (99%)	-0.05	22 (2%) 61 63	7, 22, 42, 71	8 (0%)

All (22) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	1	MET	4.8
2	B	478	HIS	4.1
2	B	454	GLY	3.4
2	B	457	LYS	3.2
2	B	144	ARG	2.9
1	A	145	LYS	2.8
2	B	456	GLU	2.8
1	A	454	GLY	2.7
2	B	322	ASN	2.6
2	B	76[A]	PHE	2.5
2	B	185	GLU	2.4
2	B	109	ARG	2.3
1	A	457	LYS	2.3
1	A	90	GLU	2.3
2	B	149	ASP	2.2
2	B	118	ASN	2.2
2	B	152	ASN	2.2
2	B	73	LYS	2.1
2	B	90	GLU	2.1
2	B	223	ALA	2.1
2	B	289	ALA	2.1
2	B	458	ALA	2.1

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

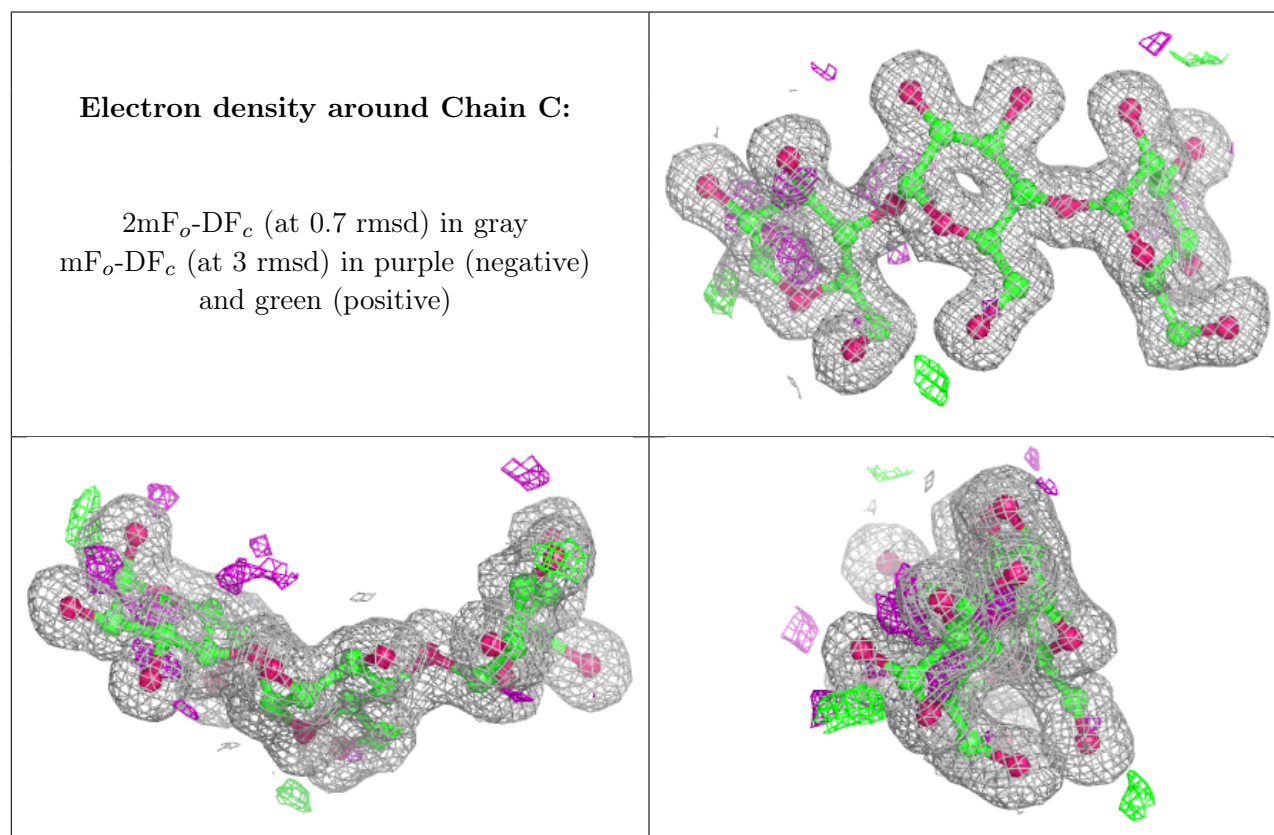
There are no non-standard protein/DNA/RNA residues in this entry.

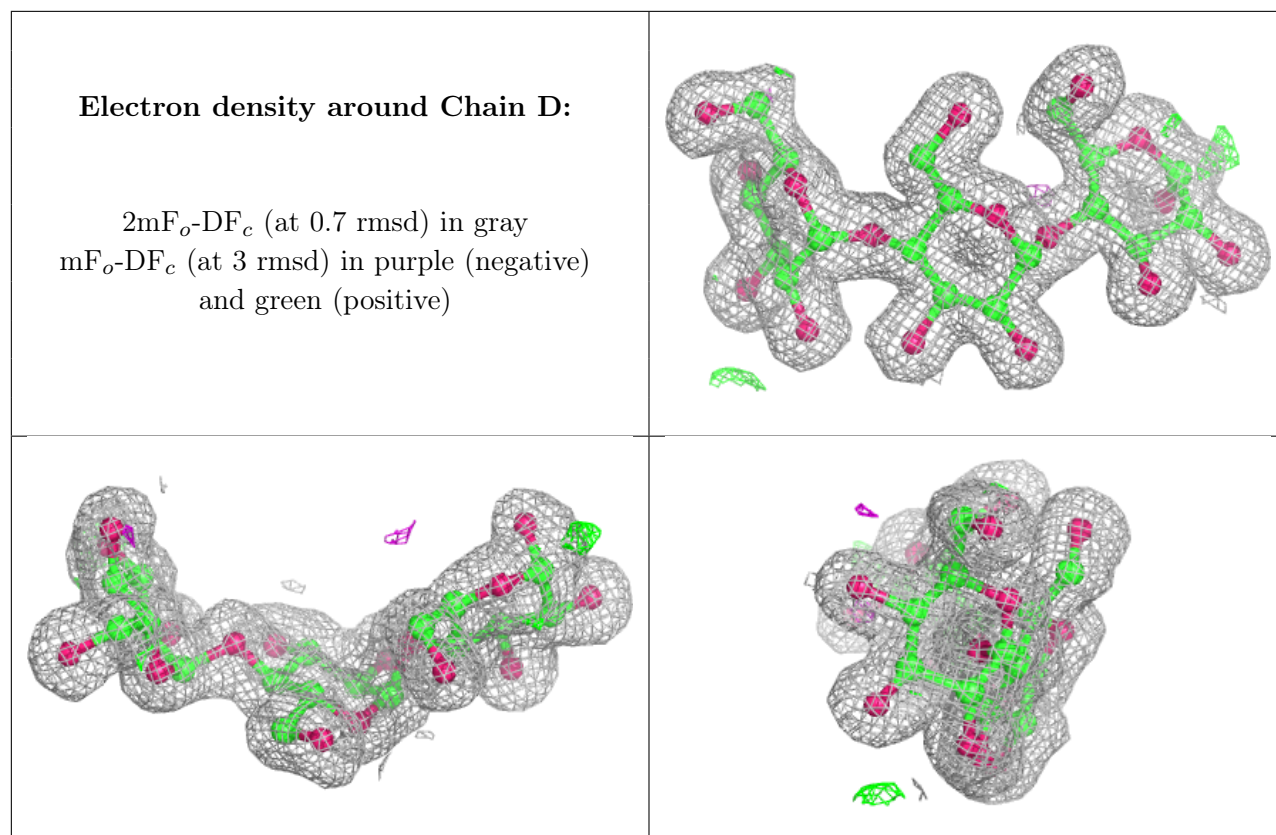
6.3 Carbohydrates [i](#)

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(\AA^2)	Q<0.9
3	GLC	C	1	12/12	0.96	0.07	12,14,21,29	0
3	GLC	D	1	12/12	0.96	0.07	15,18,26,40	0
3	GLC	D	3	11/12	0.97	0.05	17,18,21,22	0
3	GLC	C	2	11/12	0.98	0.04	8,11,12,13	0
3	GLC	D	2	11/12	0.98	0.04	13,15,16,16	0
3	GLC	C	3	11/12	0.98	0.04	11,13,16,17	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.





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(\AA^2)	Q<0.9
4	GOL	A	503	6/6	0.62	0.18	60,62,63,63	0
6	PEG	A	506	7/7	0.75	0.27	60,65,70,72	0
4	GOL	B	502	6/6	0.76	0.24	20,20,20,20	0
5	EDO	B	503	4/4	0.83	0.16	48,48,50,51	0
5	EDO	A	505	4/4	0.84	0.16	27,34,37,43	0
5	EDO	A	504	4/4	0.89	0.16	46,47,48,49	0
4	GOL	A	501	6/6	0.90	0.10	27,28,29,30	0
4	GOL	A	502	6/6	0.94	0.09	14,29,31,31	0
4	GOL	B	501	6/6	0.96	0.07	18,28,30,30	0

6.5 Other polymers [i](#)

There are no such residues in this entry.