



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 10:43 AM UTC

PDB ID : 3CKB / pdb_00003ckb
Title : B. thetaiotaomicron SusD with maltotriose
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Deposited on : 2008-03-14
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

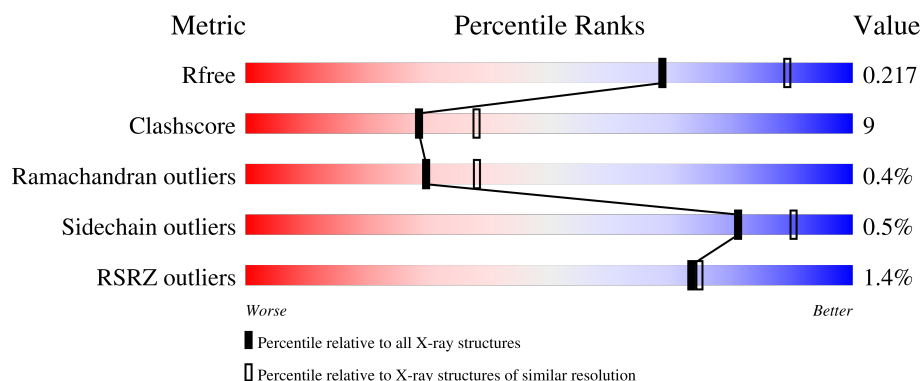
1 Overall quality at a glance

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	527	% 77% 18% • •
1	B	527	2% 70% 24% • 5%
2	C	3	33% 67%
2	D	3	100%

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 criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	GLC	C	1	X	-	-	-
2	GLC	D	1	X	-	-	-

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 8780 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 SusD.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	504	Total	C	N	O	S	0	0	0
			4049	2549	695	786	19			
1	B	501	Total	C	N	O	S	0	0	0
			4026	2534	692	781	19			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	25	GLY	-	expression tag	UNP Q8A1G2
A	100	LYS	GLU	SEE REMARK 999	UNP Q8A1G2
B	25	GLY	-	expression tag	UNP Q8A1G2
B	100	LYS	GLU	SEE REMARK 999	UNP Q8A1G2

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



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

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

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	Ca	0	0
			1	1		

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	1	Total 1	Ca 1	0	0

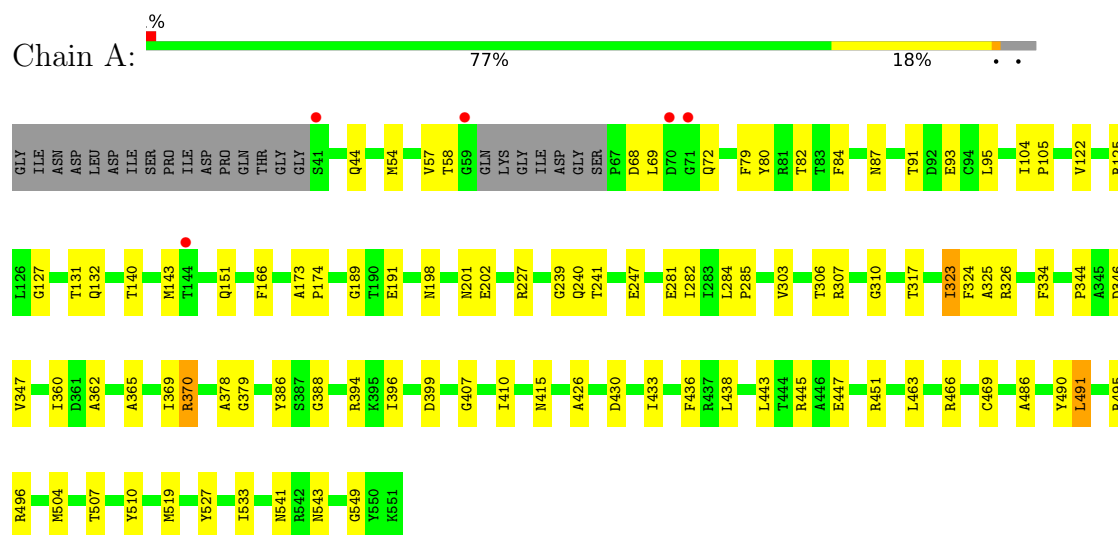
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	306	Total 306	O 306	0	0
4	B	329	Total 329	O 329	0	0

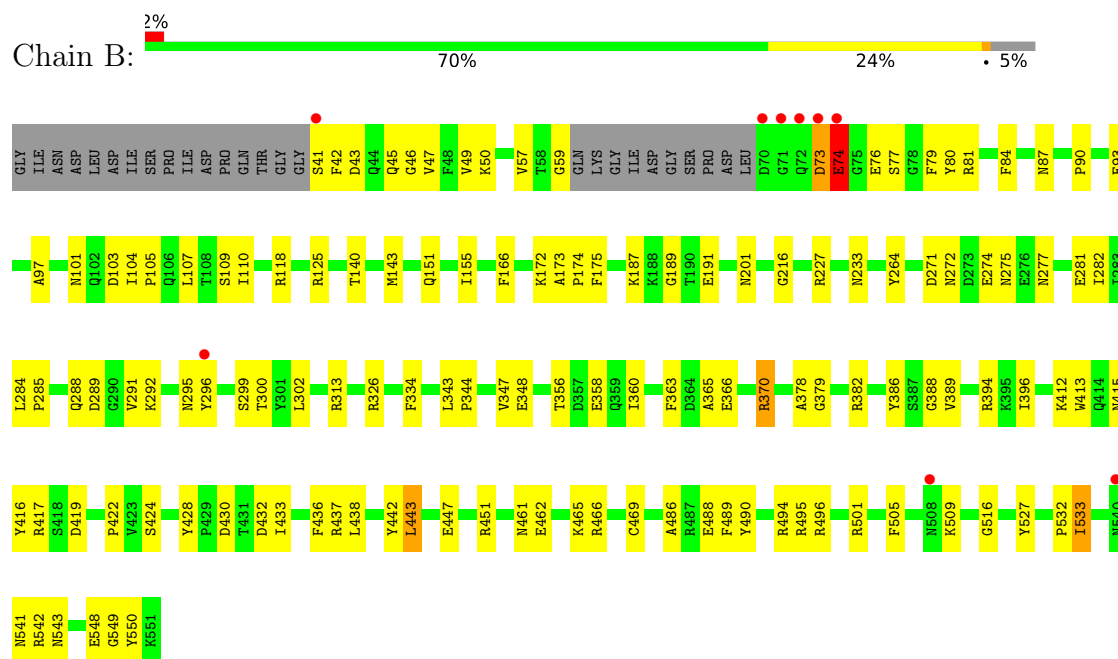
3 Residue-property plots

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

• Molecule 1: SusD



• Molecule 1: SusD



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

Chain C:  33% 67%

GLC1
GLC2
GLC3

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

Chain D:  100%

GLC1
GLC2
GLC3

4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	61.69Å 67.71Å 83.29Å 111.16° 92.94° 109.06°	Depositor
Resolution (Å)	59.50 – 2.30 59.50 – 2.30	Depositor EDS
% Data completeness (in resolution range)	91.8 (59.50-2.30) 91.7 (59.50-2.30)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.08	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.09 (at 2.29Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.180 , 0.216 0.182 , 0.217	Depositor DCC
R_{free} test set	4791 reflections (10.07%)	wwPDB-VP
Wilson B-factor (Å ²)	5.9	Xtriage
Anisotropy	0.715	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 39.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	8780	wwPDB-VP
Average B, all atoms (Å ²)	7.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.49% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: CA, GLC

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

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z > 5$	RMSZ	$\# Z > 5$
1	A	0.36	0/4142	0.86	8/5605 (0.1%)
1	B	0.37	0/4118	0.85	12/5572 (0.2%)
All	All	0.37	0/8260	0.86	20/11177 (0.2%)

There are no bond length outliers.

All (20) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	396	ILE	N-CA-C	-8.18	104.78	111.81
1	A	533	ILE	N-CA-C	-7.48	102.20	108.63
1	B	396	ILE	N-CA-C	-7.32	104.71	111.67
1	B	533	ILE	N-CA-C	-6.78	102.31	108.95
1	A	370	ARG	N-CA-C	-6.57	101.68	110.55
1	A	507	THR	N-CA-C	6.43	119.22	110.35
1	B	175	PHE	N-CA-C	6.18	120.35	109.96
1	B	505	PHE	N-CA-C	5.96	117.86	111.36
1	A	174	PRO	N-CA-C	-5.78	100.56	112.47
1	B	424	SER	N-CA-C	5.76	118.51	111.82
1	B	97	ALA	N-CA-C	5.68	117.47	111.28
1	A	323	ILE	N-CA-C	5.67	116.02	107.80
1	B	326	ARG	N-CA-C	-5.49	103.14	110.55
1	B	174	PRO	N-CA-C	-5.38	101.38	112.47
1	A	239	GLY	N-CA-C	-5.32	107.98	115.32
1	B	489	PHE	N-CA-C	5.22	118.96	112.59
1	A	91	THR	N-CA-C	5.18	117.89	109.86
1	B	77	SER	N-CA-C	-5.10	107.09	113.72
1	B	343	LEU	N-CA-C	-5.09	103.36	109.83
1	B	370	ARG	N-CA-C	-5.03	102.33	110.32

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	4049	0	3858	56	0
1	B	4026	0	3835	84	0
2	C	34	0	30	1	0
2	D	34	0	30	0	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
4	A	306	0	0	1	0
4	B	329	0	0	5	0
All	All	8780	0	7753	140	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 9.

All (140) 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:B:201:ASN:HD22	1:B:227:ARG:HH12	1.30	0.79
1:A:201:ASN:HD22	1:A:227:ARG:HH12	1.33	0.77
1:A:140:THR:HA	1:A:143:MET:HE3	1.65	0.77
1:A:69:LEU:H	1:A:72:GLN:NE2	1.84	0.75
1:A:44:GLN:HE21	1:A:151:GLN:HE22	1.35	0.73
1:B:274:GLU:CD	1:B:417:ARG:HH22	1.96	0.73
1:B:282:ILE:HG21	1:B:285:PRO:HB3	1.73	0.69
1:A:44:GLN:HE21	1:A:151:GLN:NE2	1.92	0.68
1:B:271:ASP:HB3	1:B:417:ARG:HG2	1.76	0.68
1:B:101:ASN:HB2	1:B:104:ILE:HG13	1.76	0.68
1:B:348:GLU:HB3	4:B:1234:HOH:O	1.93	0.68
1:A:95:LEU:HG	1:A:326:ARG:HG2	1.75	0.66
1:A:284:LEU:HB3	1:A:436:PHE:HB2	1.77	0.66
1:A:362:ALA:HB1	1:B:422:PRO:HG3	1.76	0.66
1:B:494:ARG:HA	4:B:1193:HOH:O	1.94	0.66
1:B:93:GLU:HG3	1:B:495:ARG:HG2	1.77	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:87:ASN:HD22	1:B:496:ARG:HB3	1.61	0.65
1:B:90:PRO:HB3	1:B:107:LEU:O	1.97	0.65
1:B:166:PHE:HB3	1:B:173:ALA:HB2	1.81	0.63
1:A:87:ASN:HD22	1:A:496:ARG:HD3	1.63	0.62
1:A:93:GLU:HG3	1:A:495:ARG:HG2	1.81	0.62
1:A:166:PHE:HB3	1:A:173:ALA:HB2	1.81	0.61
1:A:282:ILE:HG21	1:A:285:PRO:HB3	1.83	0.60
1:A:365:ALA:HB2	1:A:370:ARG:HD3	1.83	0.60
1:A:541:ASN:OD1	1:A:543:ASN:HB2	2.01	0.60
1:B:103:ASP:HB3	1:B:118:ARG:HG2	1.83	0.60
1:A:198:ASN:O	1:A:202:GLU:HG3	2.01	0.59
1:A:87:ASN:ND2	1:A:496:ARG:HD3	2.18	0.59
1:B:104:ILE:HB	1:B:105:PRO:HD3	1.83	0.58
1:A:415:ASN:HD21	1:A:430:ASP:H	1.51	0.57
1:A:191:GLU:HG3	4:A:1166:HOH:O	2.03	0.57
1:B:79:PHE:HA	1:B:125:ARG:HG2	1.87	0.56
1:B:282:ILE:CG2	1:B:285:PRO:HB3	2.35	0.56
1:B:41:SER:O	1:B:42:PHE:HB3	2.06	0.56
1:B:541:ASN:OD1	1:B:543:ASN:HB2	2.06	0.55
1:B:365:ALA:HB2	1:B:370:ARG:HD3	1.88	0.55
1:B:447:GLU:O	1:B:451:ARG:HG2	2.06	0.55
1:A:447:GLU:O	1:A:451:ARG:HG2	2.06	0.55
1:A:281:GLU:HG3	1:A:438:LEU:HB3	1.89	0.55
1:B:295:ASN:HD22	1:B:300:THR:HG21	1.72	0.54
1:B:344:PRO:HB2	1:B:347:VAL:HG23	1.89	0.53
1:A:201:ASN:HD22	1:A:227:ARG:NH1	2.05	0.53
1:A:415:ASN:ND2	1:A:430:ASP:H	2.05	0.53
1:B:291:VAL:HG23	1:B:292:LYS:HG3	1.90	0.52
1:B:379:GLY:O	1:B:469:CYS:HA	2.09	0.52
1:B:415:ASN:ND2	1:B:430:ASP:H	2.08	0.52
1:B:442:TYR:CE1	1:B:462:GLU:HG2	2.45	0.52
1:B:81:ARG:HG3	4:B:1239:HOH:O	2.09	0.51
1:B:288:GLN:HG3	1:B:299:SER:OG	2.10	0.51
1:B:140:THR:HA	1:B:143:MET:HE3	1.91	0.51
1:B:284:LEU:HB3	1:B:436:PHE:HB2	1.91	0.51
1:B:415:ASN:HD21	1:B:430:ASP:H	1.58	0.51
1:B:281:GLU:HG3	1:B:438:LEU:HB3	1.92	0.50
1:B:187:LYS:HG3	1:B:191:GLU:HG2	1.94	0.50
1:B:358:GLU:CD	1:B:358:GLU:H	2.19	0.50
1:B:101:ASN:HB2	1:B:104:ILE:CG1	2.41	0.50
1:A:87:ASN:ND2	1:A:496:ARG:HB3	2.27	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:486:ALA:O	1:A:490:TYR:HB3	2.11	0.49
1:B:437:ARG:NH2	1:B:488:GLU:OE2	2.38	0.49
1:B:45:GLN:O	1:B:49:VAL:HG23	2.12	0.49
1:A:310:GLY:O	1:A:360:ILE:HD13	2.13	0.49
1:B:59:GLY:N	4:B:1330:HOH:O	2.45	0.49
1:B:109:SER:O	1:B:516:GLY:HA2	2.12	0.49
1:A:87:ASN:HD22	1:A:496:ARG:HB3	1.78	0.48
1:A:344:PRO:HB2	1:A:347:VAL:HG23	1.95	0.48
1:B:80:TYR:CE2	1:B:433:ILE:HG21	2.49	0.48
1:B:334:PHE:CZ	1:B:378:ALA:HB2	2.49	0.48
1:B:87:ASN:ND2	1:B:496:ARG:HB3	2.28	0.47
1:A:247:GLU:HG3	1:A:445:ARG:HG3	1.95	0.47
1:B:172:LYS:HD2	1:B:533:ILE:HD11	1.95	0.47
1:B:187:LYS:CG	1:B:191:GLU:HG2	2.45	0.47
1:A:369:ILE:HD12	1:A:388:GLY:HA2	1.97	0.47
1:B:386:TYR:CZ	1:B:388:GLY:HA3	2.49	0.47
1:B:486:ALA:O	1:B:490:TYR:HB3	2.15	0.47
1:B:527:TYR:HB2	1:B:548:GLU:O	2.14	0.47
1:B:295:ASN:HB2	1:B:428:TYR:CE2	2.50	0.47
1:B:542:ARG:HG3	1:B:542:ARG:HH11	1.79	0.47
1:B:110:ILE:HG22	1:B:110:ILE:O	2.15	0.46
1:A:57:VAL:HG22	1:A:58:THR:N	2.31	0.46
1:B:417:ARG:HD3	1:B:419:ASP:OD1	2.15	0.46
1:A:443:LEU:HG	1:A:463:LEU:CD1	2.45	0.46
1:B:87:ASN:HD22	1:B:496:ARG:HD3	1.81	0.46
1:A:68:ASP:HA	1:A:72:GLN:HE22	1.81	0.46
1:A:324:PHE:CD1	1:A:407:GLY:HA3	2.51	0.45
1:B:87:ASN:HD21	1:B:496:ARG:HH11	1.65	0.45
1:B:233:ASN:OD1	1:B:501:ARG:HD3	2.16	0.45
1:B:73:ASP:CG	1:B:74:GLU:H	2.22	0.45
1:B:363:PHE:CE1	1:B:389:VAL:HG11	2.52	0.45
1:A:527:TYR:CG	1:A:549:GLY:HA3	2.51	0.45
1:B:275:ASN:C	1:B:275:ASN:ND2	2.75	0.45
1:B:101:ASN:N	1:B:101:ASN:HD22	2.14	0.45
1:B:313:ARG:HB3	1:B:416:TYR:CE2	2.51	0.45
1:A:466:ARG:HD2	1:A:466:ARG:C	2.42	0.44
1:B:275:ASN:C	1:B:275:ASN:HD22	2.24	0.44
1:A:346:ASP:OD2	1:A:346:ASP:C	2.60	0.44
1:B:289:ASP:OD1	1:B:291:VAL:HG22	2.17	0.44
1:A:306:THR:HG22	1:A:323:ILE:HB	1.98	0.44
1:B:43:ASP:O	1:B:47:VAL:HG23	2.16	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:73:ASP:HB3	1:B:76:GLU:CD	2.42	0.44
2:C:1:GLC:H62	2:C:2:GLC:H5	2.00	0.44
1:A:386:TYR:CZ	1:A:388:GLY:HA3	2.53	0.44
1:B:46:GLY:O	1:B:50:LYS:HG3	2.17	0.44
1:B:57:VAL:HG23	1:B:296:TYR:O	2.18	0.44
1:B:264:TYR:CZ	1:B:382:ARG:HG2	2.53	0.44
1:A:240:GLN:NE2	1:A:241:THR:O	2.50	0.44
1:A:82:THR:HB	1:A:122:VAL:HB	1.99	0.44
1:B:275:ASN:ND2	1:B:277:ASN:H	2.16	0.43
1:A:394:ARG:HG2	1:A:394:ARG:HH11	1.82	0.43
1:A:317:THR:HG23	1:A:426:ALA:O	2.19	0.43
1:A:504:MET:HB3	1:A:510:TYR:HB3	2.00	0.43
1:B:356:THR:O	1:B:360:ILE:HG13	2.18	0.43
1:A:127:GLY:O	1:A:131:THR:HG23	2.19	0.43
1:A:527:TYR:CD1	1:A:527:TYR:C	2.96	0.43
1:B:527:TYR:CG	1:B:549:GLY:HA3	2.53	0.43
1:A:325:ALA:HB2	1:A:410:ILE:HD11	2.01	0.43
1:B:302:LEU:HD13	1:B:413:TRP:CE2	2.54	0.43
1:A:303:VAL:O	1:A:307:ARG:HG2	2.18	0.43
1:B:216:GLY:HA2	1:B:285:PRO:HG2	2.01	0.43
1:A:84:PHE:HZ	1:A:491:LEU:HD13	1.84	0.42
1:B:412:LYS:HD2	1:B:490:TYR:CZ	2.54	0.42
1:A:334:PHE:CZ	1:A:378:ALA:HB2	2.54	0.42
1:A:54:MET:HE1	1:A:132:GLN:HB2	2.01	0.42
1:A:399:ASP:HA	1:A:519:MET:HG3	2.01	0.42
1:B:84:PHE:C	1:B:84:PHE:CD1	2.97	0.42
1:A:282:ILE:CG2	1:A:285:PRO:HB3	2.49	0.42
1:B:532:PRO:HB3	1:B:550:TYR:CD2	2.55	0.42
1:B:151:GLN:O	1:B:155:ILE:HG12	2.20	0.42
1:A:104:ILE:HB	1:A:105:PRO:HD3	2.01	0.42
1:B:103:ASP:HB3	1:B:118:ARG:CG	2.48	0.42
1:B:264:TYR:CE2	1:B:382:ARG:HG2	2.55	0.42
1:A:79:PHE:HA	1:A:125:ARG:HG2	2.02	0.41
1:B:201:ASN:HD22	1:B:227:ARG:NH1	2.07	0.41
1:A:80:TYR:CE2	1:A:433:ILE:HG21	2.55	0.41
1:A:379:GLY:O	1:A:469:CYS:HA	2.20	0.41
1:B:394:ARG:HH11	1:B:394:ARG:HG3	1.86	0.41
1:B:443:LEU:HD12	1:B:443:LEU:HA	1.87	0.40
1:B:466:ARG:HD2	1:B:466:ARG:C	2.46	0.40
1:B:461:ASN:O	1:B:465:LYS:HG3	2.22	0.40
1:B:272:ASN:OD1	1:B:432:ASP:HB2	2.21	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:509:LYS:HD2	4:B:1412:HOH:O	2.22	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	500/527 (95%)	478 (96%)	21 (4%)	1 (0%)	43	55
1	B	497/527 (94%)	472 (95%)	22 (4%)	3 (1%)	21	27
All	All	997/1054 (95%)	950 (95%)	43 (4%)	4 (0%)	30	38

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	74	GLU
1	A	189	GLY
1	B	73	ASP
1	B	189	GLY

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	423/441 (96%)	422 (100%)	1 (0%)	87	94

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	B	420/441 (95%)	417 (99%)	3 (1%)	76	87
All	All	843/882 (96%)	839 (100%)	4 (0%)	81	90

All (4) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	491	LEU
1	B	74	GLU
1	B	366	GLU
1	B	443	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (23) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	44	GLN
1	A	72	GLN
1	A	87	ASN
1	A	102	GLN
1	A	151	GLN
1	A	198	ASN
1	A	201	ASN
1	A	415	ASN
1	A	540	ASN
1	A	543	ASN
1	B	44	GLN
1	B	87	ASN
1	B	99	GLN
1	B	102	GLN
1	B	151	GLN
1	B	198	ASN
1	B	201	ASN
1	B	275	ASN
1	B	295	ASN
1	B	331	GLN
1	B	415	ASN
1	B	540	ASN
1	B	543	ASN

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$
2	GLC	C	1	2	12,12,12	0.93	0	17,17,17	1.05	2 (11%)
2	GLC	C	2	2	11,11,12	1.23	1 (9%)	15,15,17	0.81	0
2	GLC	C	3	2	11,11,12	1.14	0	15,15,17	0.81	0
2	GLC	D	1	2	12,12,12	1.04	0	17,17,17	0.96	1 (5%)
2	GLC	D	2	2	11,11,12	1.30	1 (9%)	15,15,17	0.84	0
2	GLC	D	3	2	11,11,12	1.16	1 (9%)	15,15,17	0.80	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	GLC	C	1	2	1/1/5/5	0/2/22/22	0/1/1/1
2	GLC	C	2	2	-	0/2/19/22	0/1/1/1
2	GLC	C	3	2	-	2/2/19/22	0/1/1/1
2	GLC	D	1	2	1/1/5/5	0/2/22/22	0/1/1/1
2	GLC	D	2	2	-	2/2/19/22	0/1/1/1
2	GLC	D	3	2	-	2/2/19/22	0/1/1/1

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	2	GLC	O5-C1	2.23	1.47	1.43
2	C	2	GLC	O5-C1	2.11	1.47	1.43
2	D	3	GLC	O5-C1	2.00	1.47	1.43

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	1	GLC	C1-O5-C5	-2.15	109.48	113.65
2	D	1	GLC	C6-C5-C4	-2.04	108.02	113.02
2	C	1	GLC	C1-C2-C3	-2.02	106.23	110.36

All (2) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
2	C	1	GLC	C1
2	D	1	GLC	C1

All (6) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	D	3	GLC	O5-C5-C6-O6
2	D	2	GLC	C4-C5-C6-O6
2	D	3	GLC	C4-C5-C6-O6
2	D	2	GLC	O5-C5-C6-O6
2	C	3	GLC	C4-C5-C6-O6
2	C	3	GLC	O5-C5-C6-O6

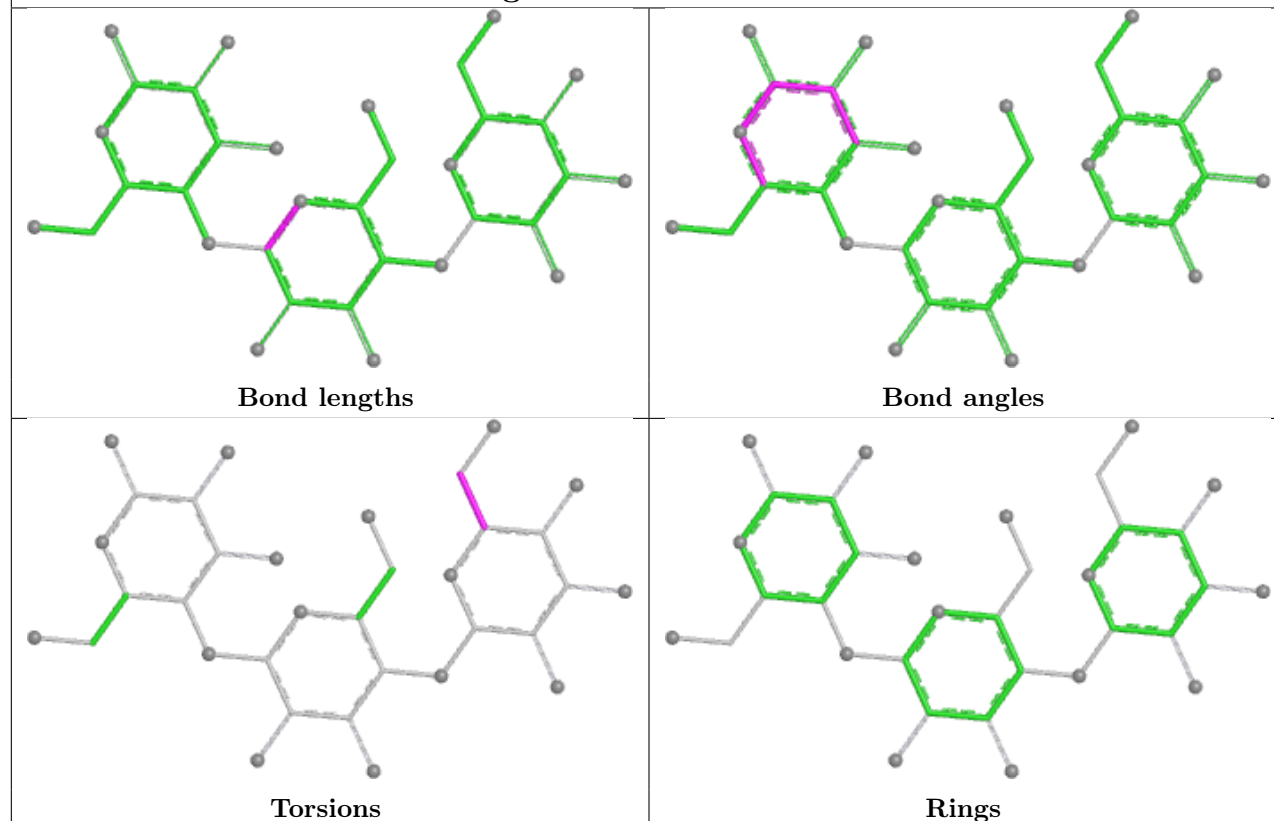
There are no ring outliers.

2 monomers are involved in 1 short contact:

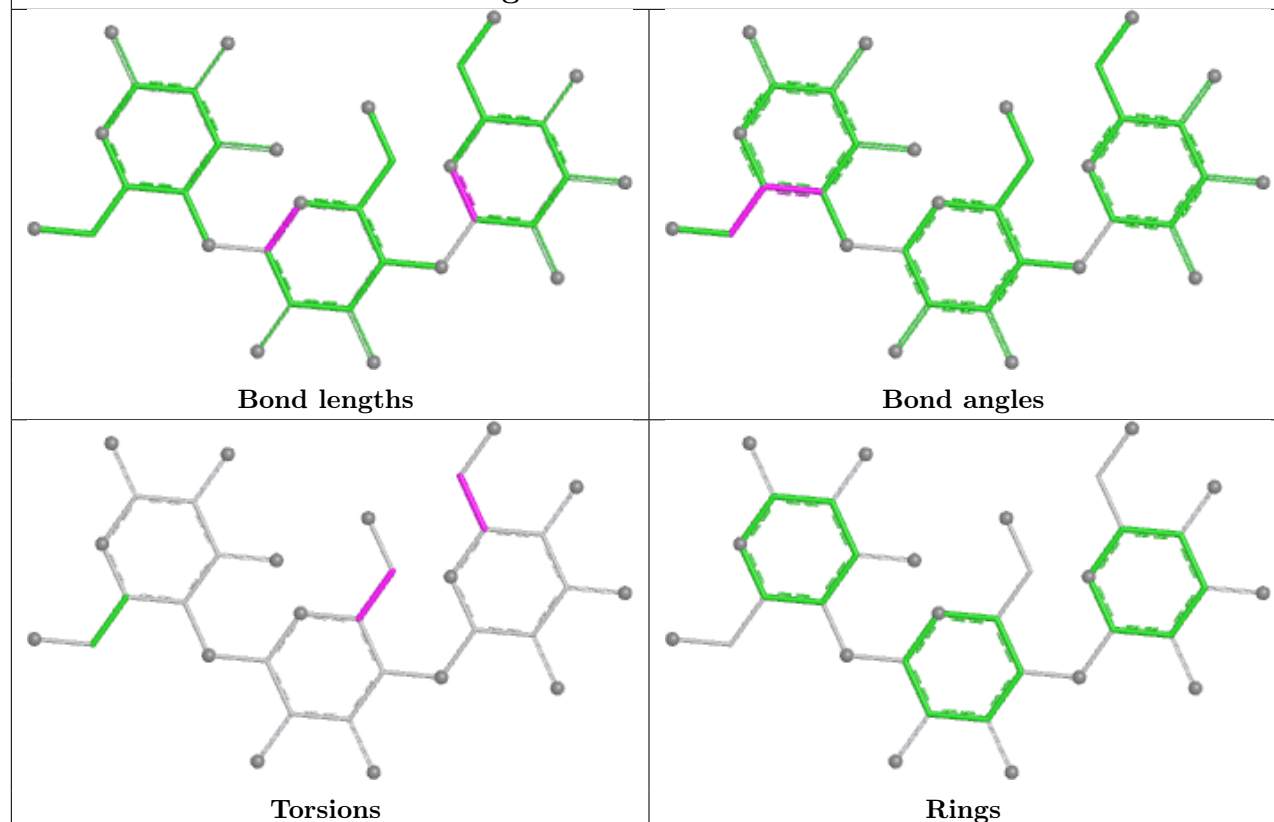
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	2	GLC	1	0
2	C	1	GLC	1	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for oligosaccharide.

Oligosaccharide Chain C



Oligosaccharide Chain D



5.6 Ligand geometry

Of 2 ligands modelled in this entry, 2 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

There are no chain breaks in this entry.

6 Fit of model and data [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	504/527 (95%)	-0.36	5 (0%) 79 80	1, 7, 18, 29	0
1	B	501/527 (95%)	-0.35	9 (1%) 67 69	1, 6, 17, 33	0
All	All	1005/1054 (95%)	-0.36	14 (1%) 73 75	1, 6, 17, 33	0

All (14) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	70	ASP	3.4
1	B	71	GLY	3.1
1	B	73	ASP	3.0
1	B	41	SER	2.9
1	B	508	ASN	2.6
1	A	41	SER	2.5
1	A	71	GLY	2.3
1	A	70	ASP	2.3
1	B	72	GLN	2.2
1	A	59	GLY	2.2
1	B	74	GLU	2.1
1	B	540	ASN	2.1
1	A	144	THR	2.1
1	B	296	TYR	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.

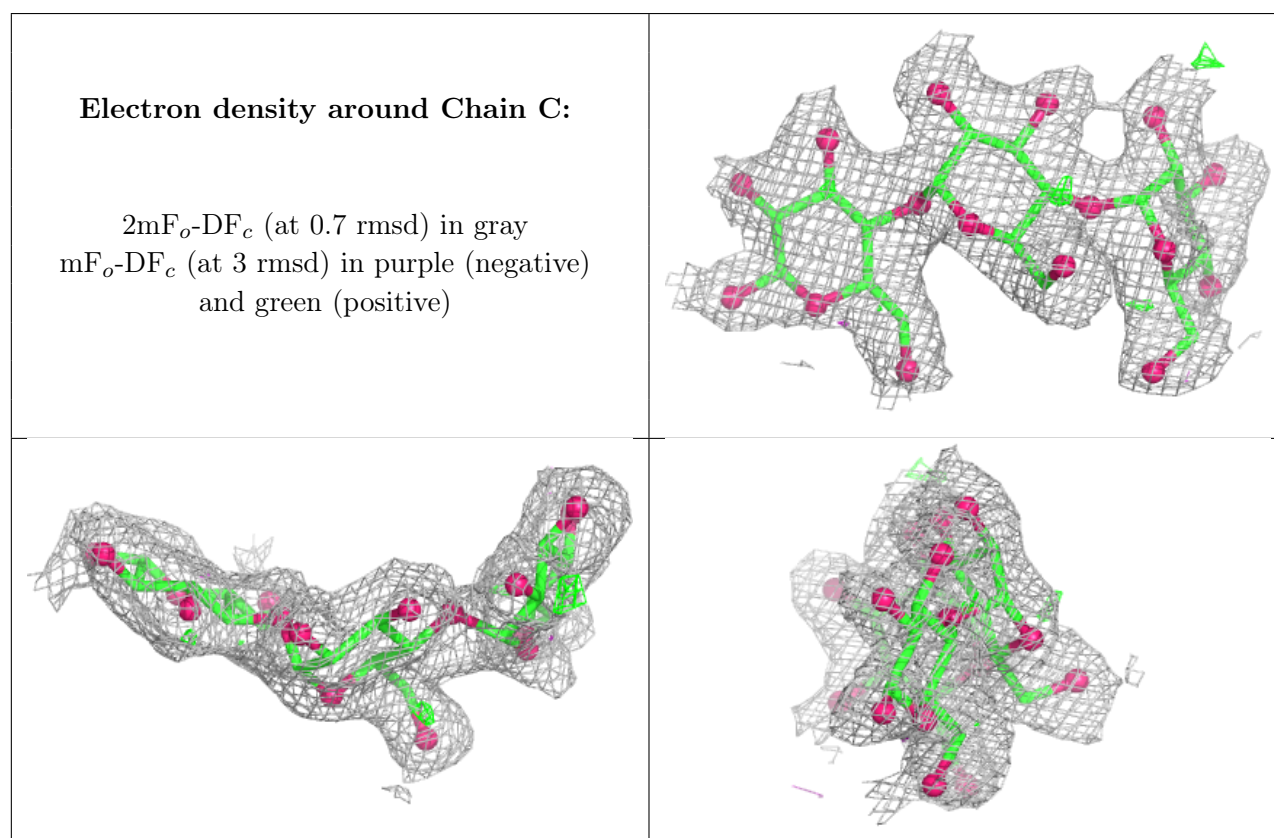
6.3 Carbohydrates [i](#)

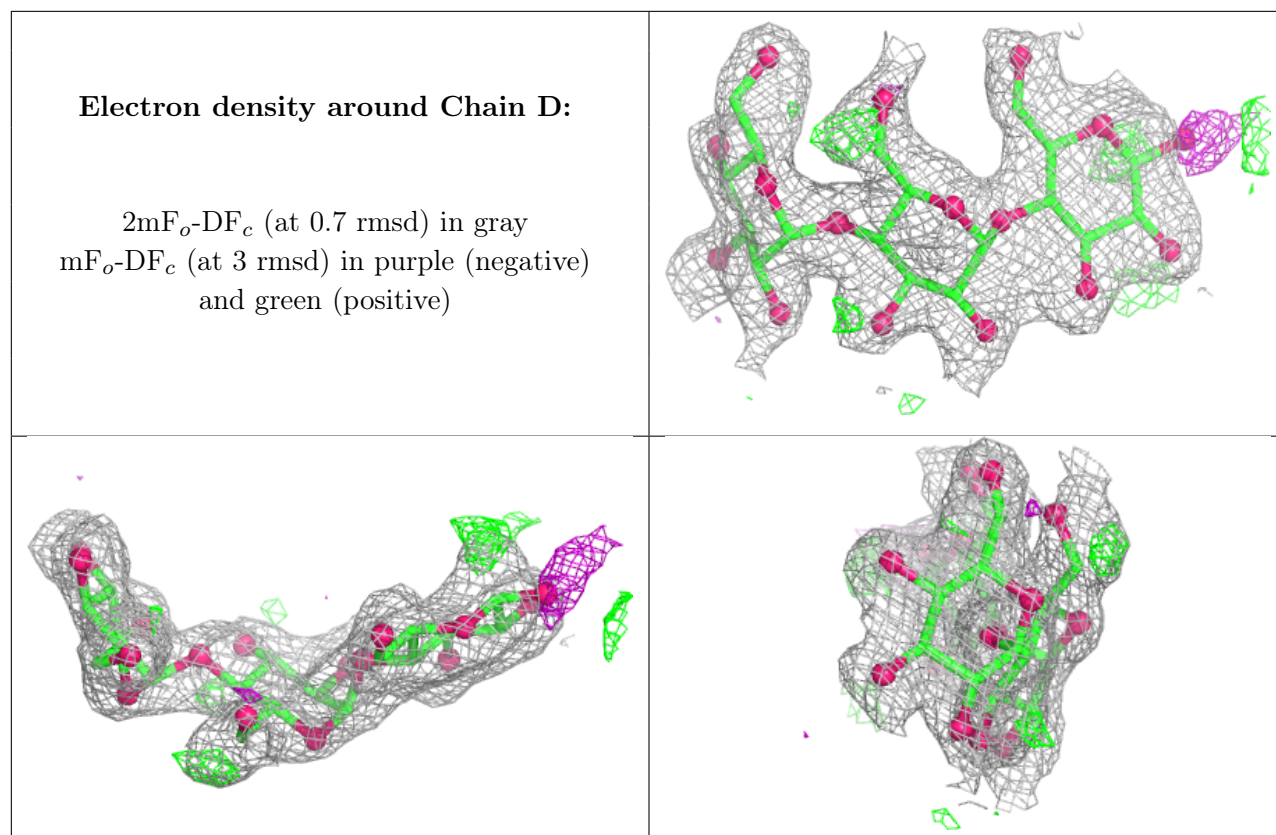
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum,

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(Å ²)	Q<0.9
2	GLC	D	3	11/12	0.82	0.12	21,23,25,25	0
2	GLC	C	1	12/12	0.83	0.12	17,18,21,21	0
2	GLC	D	2	11/12	0.85	0.11	18,19,21,25	0
2	GLC	D	1	12/12	0.86	0.11	15,18,20,27	0
2	GLC	C	3	11/12	0.87	0.10	14,15,18,22	0
2	GLC	C	2	11/12	0.94	0.08	10,13,14,16	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($\text{\AA}^2$)	Q<0.9
3	CA	A	600	1/1	0.98	0.02	4,4,4,4	0
3	CA	B	700	1/1	0.99	0.02	4,4,4,4	0

6.5 Other polymers [i](#)

There are no such residues in this entry.