



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 10, 2026 – 12:56 AM UTC

PDB ID : 4EN8 / pdb_00004en8
Title : Crystal structure of HA70 (HA3) subcomponent of Clostridium botulinum type C progenitor toxin in complex with alpha 2-6-sialyllactose
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Deposited on : 2012-04-12
Resolution : 2.60 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 Phenix2.0
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

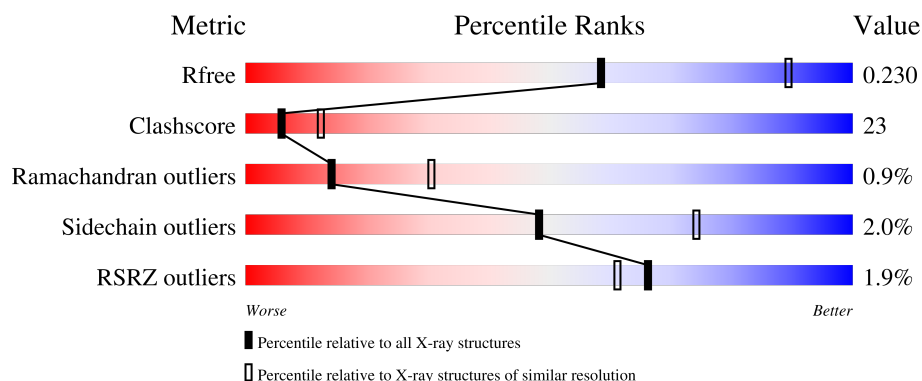
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.60 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}	180053	4008 (2.60-2.60)
Clashscore	190562	4347 (2.60-2.60)
Ramachandran outliers	187476	4277 (2.60-2.60)
Sidechain outliers	187428	4277 (2.60-2.60)
RSRZ outliers	180081	4008 (2.60-2.60)

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	224	 2% 47% 27% 24%
2	B	420	 2% 66% 30%
3	C	3	 67% 33%

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
4	MRD	A	301	-	-	X	-
4	MRD	A	303	-	-	X	-

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 5161 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 Hemagglutinin components HA-22/23/53.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	170	Total	C	N	O	S	0	0	0
			1382	882	226	271	3			

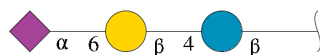
There are 21 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-20	ILE	-	expression tag	UNP P46085
A	-19	SER	-	expression tag	UNP P46085
A	-18	GLU	-	expression tag	UNP P46085
A	-17	PHE	-	expression tag	UNP P46085
A	-16	ASP	-	expression tag	UNP P46085
A	-15	TYR	-	expression tag	UNP P46085
A	-14	LYS	-	expression tag	UNP P46085
A	-13	ASP	-	expression tag	UNP P46085
A	-12	HIS	-	expression tag	UNP P46085
A	-11	ASP	-	expression tag	UNP P46085
A	-10	ILE	-	expression tag	UNP P46085
A	-9	ASP	-	expression tag	UNP P46085
A	-8	TYR	-	expression tag	UNP P46085
A	-7	LYS	-	expression tag	UNP P46085
A	-6	ASP	-	expression tag	UNP P46085
A	-5	ASP	-	expression tag	UNP P46085
A	-4	ASP	-	expression tag	UNP P46085
A	-3	ASP	-	expression tag	UNP P46085
A	-2	LYS	-	expression tag	UNP P46085
A	-1	TRP	-	expression tag	UNP P46085
A	0	ILE	-	expression tag	UNP P46085

- Molecule 2 is a protein called Hemagglutinin components HA-22/23/53.

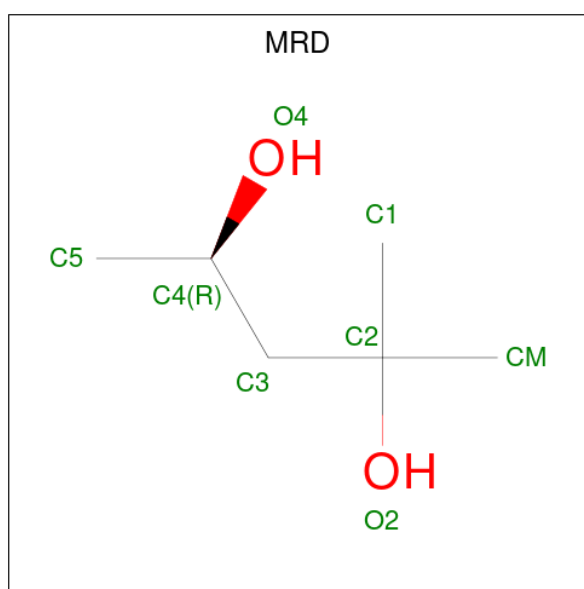
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	420	Total	C	N	O	S	0	0	0
			3339	2109	553	672	5			

- Molecule 3 is an oligosaccharide called N-acetyl-alpha-neuraminic acid-(2-6)-beta-D-galactopyranose-(1-4)-beta-D-glucopyranose.



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
3	C	3	Total	C	N	O	0	0	0
			43	23	1	19			

- Molecule 4 is (4R)-2-METHYLPENTANE-2,4-DIOL (CCD ID: MRD) (formula: C₆H₁₄O₂).

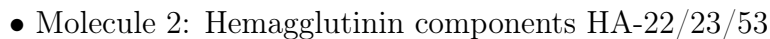


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			8	6	2		
4	A	1	Total	C	O	0	0
			8	6	2		
4	A	1	Total	C	O	0	0
			8	6	2		
4	B	1	Total	C	O	0	0
			8	6	2		
4	B	1	Total	C	O	0	0
			8	6	2		

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	94	Total 94	O 94	0	0
5	B	263	Total 263	O 263	0	0

- Molecule 1: Hemagglutinin components HA-22/23/53





4 Data and refinement statistics

Property	Value	Source
Space group	P 63	Depositor
Cell constants a, b, c, α , β , γ	177.28Å 177.28Å 80.79Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	34.67 – 2.60 34.67 – 2.60	Depositor EDS
% Data completeness (in resolution range)	99.7 (34.67-2.60) 99.8 (34.67-2.60)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.47 (at 2.61Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.193 , 0.235 0.193 , 0.230	Depositor DCC
R_{free} test set	4505 reflections (10.10%)	wwPDB-VP
Wilson B-factor (Å ²)	49.6	Xtriage
Anisotropy	0.066	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 48.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.42$, $\langle L^2 \rangle = 0.25$	Xtriage
Estimated twinning fraction	0.043 for h,-h-k,-l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	5161	wwPDB-VP
Average B, all atoms (Å ²)	52.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 11.76% 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: GAL, BGC, SIA, MRD

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	0.42	0/1411	0.97	4/1914 (0.2%)
2	B	0.40	0/3405	0.96	18/4642 (0.4%)
All	All	0.40	0/4816	0.97	22/6556 (0.3%)

There are no bond length outliers.

All (22) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	501	ILE	N-CA-C	-7.19	103.65	113.00
2	B	249	LEU	N-CA-C	-7.17	104.50	113.18
2	B	219	GLY	N-CA-C	6.72	118.90	112.04
1	A	50	ASN	N-CA-C	6.71	120.63	111.17
2	B	461	SER	N-CA-C	6.66	118.20	111.07
2	B	308	ASN	N-CA-C	-6.29	102.06	110.55
2	B	245	ILE	N-CA-C	6.00	117.41	108.46
2	B	465	SER	N-CA-C	5.98	116.31	107.88
2	B	557	ARG	N-CA-C	5.59	119.87	112.72
2	B	452	ILE	N-CA-C	5.46	116.60	108.46
2	B	540	ILE	N-CA-C	5.42	114.86	108.96
2	B	261	ASN	CA-C-N	5.37	125.03	119.56
2	B	261	ASN	C-N-CA	5.37	125.03	119.56
2	B	435	SER	N-CA-C	5.18	118.00	109.76
1	A	153	PHE	N-CA-C	5.18	115.37	108.38
2	B	311	GLU	N-CA-C	-5.09	100.94	109.24
1	A	41	GLN	N-CA-C	5.08	118.79	112.59
1	A	165	PHE	N-CA-C	5.01	117.23	108.76
2	B	207	LEU	CA-C-N	5.01	126.10	119.84
2	B	207	LEU	C-N-CA	5.01	126.10	119.84
2	B	246	GLN	CA-C-N	5.00	125.44	120.04
2	B	246	GLN	C-N-CA	5.00	125.44	120.04

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	1382	0	1336	69	0
2	B	3339	0	3275	141	0
3	C	43	0	37	6	0
4	A	24	0	42	16	0
4	B	16	0	28	4	0
5	A	94	0	0	20	0
5	B	263	0	0	24	0
All	All	5161	0	4718	218	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 23.

All (218) 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:161:LYS:HD3	4:A:303:MRD:HMC2	1.56	0.85
2:B:512:LEU:HD22	2:B:538:LEU:HD21	1.55	0.85
2:B:216:ILE:HD12	2:B:369:ILE:HG12	1.59	0.83
2:B:378:ASN:H	2:B:378:ASN:HD22	1.27	0.83
2:B:336:LYS:HE2	2:B:362:THR:CG2	2.09	0.82
2:B:545:ASN:HD21	2:B:548:ALA:HB3	1.43	0.82
2:B:545:ASN:HD21	2:B:548:ALA:CB	1.95	0.79
2:B:221:GLY:O	2:B:362:THR:HB	1.84	0.76
2:B:250:ARG:HA	2:B:250:ARG:NH1	2.01	0.76
1:A:85:GLN:NE2	1:A:92:THR:H	1.84	0.75
2:B:555:ARG:HD3	5:B:815:HOH:O	1.85	0.75
1:A:130:LEU:O	1:A:131:GLN:HB2	1.87	0.74
2:B:336:LYS:HE2	2:B:362:THR:HG23	1.71	0.73
2:B:373:ILE:HD12	2:B:452:ILE:HD13	1.70	0.72
2:B:512:LEU:HD22	2:B:538:LEU:CD2	2.19	0.71
2:B:261:ASN:CG	2:B:264:ILE:HG22	2.16	0.71

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:273:ARG:NH1	2:B:276:ASP:OD2	2.24	0.71
2:B:261:ASN:OD1	2:B:264:ILE:HG22	1.90	0.71
1:A:37:SER:O	1:A:38:ARG:HG2	1.92	0.70
2:B:383:SER:HB3	5:B:1015:HOH:O	1.91	0.69
2:B:547:ILE:HG13	2:B:548:ALA:N	2.08	0.69
2:B:545:ASN:O	2:B:545:ASN:ND2	2.26	0.68
4:A:301:MRD:H5C2	5:A:472:HOH:O	1.92	0.68
1:A:15:SER:HA	5:A:430:HOH:O	1.94	0.68
1:A:92:THR:HA	5:A:466:HOH:O	1.94	0.67
2:B:567:ARG:HD3	2:B:601:TYR:CE1	2.30	0.67
2:B:612:ASN:ND2	2:B:614:ASN:H	1.93	0.67
1:A:55:ILE:HD13	4:A:301:MRD:H5C3	1.77	0.67
2:B:428:ILE:HG23	2:B:455:ILE:HB	1.76	0.67
1:A:15:SER:N	1:A:18:ASN:ND2	2.44	0.66
2:B:264:ILE:HB	5:B:905:HOH:O	1.95	0.66
2:B:336:LYS:HE2	2:B:362:THR:HG21	1.78	0.66
1:A:128:LYS:HA	5:A:442:HOH:O	1.95	0.65
1:A:88:PHE:HB3	5:A:407:HOH:O	1.96	0.65
2:B:512:LEU:CD2	2:B:538:LEU:HD21	2.27	0.65
1:A:108:PRO:HB2	1:A:111:LYS:HE2	1.79	0.65
5:A:406:HOH:O	4:B:705:MRD:H1C1	1.97	0.64
2:B:545:ASN:ND2	2:B:548:ALA:HB3	2.12	0.64
1:A:47:ILE:HD13	2:B:362:THR:HG22	1.79	0.64
2:B:460:ARG:CZ	3:C:1:BGC:H3	2.29	0.63
2:B:294:GLU:HB3	2:B:304:MET:HE3	1.81	0.63
2:B:435:SER:HB3	2:B:475:LYS:HG2	1.80	0.63
2:B:226:ASN:HD21	2:B:355:ASP:H	1.45	0.63
2:B:291:LYS:HD3	5:B:924:HOH:O	2.00	0.62
1:A:108:PRO:HG2	1:A:111:LYS:HB2	1.80	0.62
2:B:597:ASN:HB3	5:B:846:HOH:O	2.00	0.62
2:B:515:LEU:HD11	2:B:538:LEU:HD22	1.82	0.61
1:A:168:GLN:HB2	5:A:420:HOH:O	2.01	0.61
2:B:545:ASN:O	2:B:546:ASN:C	2.44	0.61
2:B:547:ILE:HG13	2:B:548:ALA:H	1.66	0.60
2:B:612:ASN:HD22	2:B:614:ASN:H	1.48	0.60
1:A:66:THR:HG23	5:A:433:HOH:O	2.00	0.60
1:A:85:GLN:HE22	1:A:92:THR:H	1.49	0.59
2:B:505:ALA:HB3	5:B:918:HOH:O	2.01	0.59
2:B:598:GLY:N	5:B:846:HOH:O	2.35	0.59
2:B:250:ARG:HA	2:B:250:ARG:CZ	2.32	0.59
2:B:528:PRO:CG	2:B:534:ILE:HB	2.34	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:264:ILE:HG23	2:B:265:PHE:N	2.18	0.58
2:B:545:ASN:HA	5:B:1023:HOH:O	2.03	0.58
2:B:552:PHE:CD2	2:B:595:LEU:HD12	2.39	0.58
1:A:58:ASP:OD2	1:A:59:LEU:N	2.29	0.57
1:A:59:LEU:HD11	4:A:302:MRD:HMC3	1.86	0.57
2:B:528:PRO:HG2	2:B:534:ILE:HB	1.86	0.57
1:A:111:LYS:HD2	5:A:452:HOH:O	2.04	0.57
1:A:131:GLN:C	5:A:453:HOH:O	2.46	0.57
2:B:216:ILE:CD1	2:B:369:ILE:HG23	2.35	0.57
4:A:303:MRD:HMC1	4:A:303:MRD:H5C3	1.86	0.56
1:A:85:GLN:HE22	1:A:92:THR:CB	2.18	0.56
2:B:595:LEU:HB3	2:B:601:TYR:CE2	2.41	0.56
1:A:30:ARG:NH1	1:A:175:GLU:OE2	2.39	0.55
1:A:161:LYS:HD3	4:A:303:MRD:CM	2.31	0.55
2:B:460:ARG:NH1	3:C:1:BGC:H3	2.21	0.55
1:A:40:ASN:HD22	1:A:40:ASN:H	1.55	0.55
2:B:314:GLU:OE2	2:B:326:LYS:HE3	2.07	0.55
2:B:545:ASN:O	2:B:547:ILE:N	2.40	0.55
1:A:69:TYR:CE1	1:A:140:LEU:HD13	2.42	0.54
1:A:130:LEU:N	5:A:474:HOH:O	2.38	0.54
2:B:361:GLN:NE2	5:B:807:HOH:O	2.41	0.54
1:A:102:GLU:HB2	4:A:301:MRD:HMC2	1.90	0.54
1:A:59:LEU:HD11	4:A:302:MRD:CM	2.36	0.54
2:B:305:LYS:HE2	5:B:928:HOH:O	2.07	0.54
2:B:557:ARG:HD2	5:B:926:HOH:O	2.08	0.54
2:B:425:PRO:HG3	2:B:460:ARG:O	2.07	0.53
2:B:515:LEU:CD1	2:B:538:LEU:HD22	2.39	0.53
3:C:1:BGC:H4	3:C:2:GAL:O2	2.08	0.53
1:A:131:GLN:O	1:A:132:ASN:HB2	2.09	0.53
4:A:301:MRD:H1C1	5:A:454:HOH:O	2.08	0.52
1:A:69:TYR:CD1	1:A:140:LEU:HD13	2.44	0.52
2:B:591:GLU:HG2	5:B:1022:HOH:O	2.08	0.52
1:A:50:ASN:HB3	5:A:406:HOH:O	2.08	0.52
1:A:67:PRO:HD3	5:A:486:HOH:O	2.08	0.52
1:A:37:SER:C	1:A:38:ARG:HG2	2.34	0.52
1:A:59:LEU:HD23	1:A:99:ILE:O	2.09	0.52
2:B:256:LEU:HB2	2:B:330:TYR:HB2	1.92	0.52
2:B:560:GLN:HB3	2:B:609:GLU:HB3	1.92	0.52
2:B:512:LEU:HD11	2:B:536:TYR:HB3	1.92	0.52
1:A:17:ASN:O	1:A:133:THR:HA	2.10	0.52
1:A:46:ASN:ND2	1:A:54:ALA:HB1	2.25	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:551:ILE:HB	2:B:622:ASN:HB2	1.92	0.51
2:B:378:ASN:H	2:B:378:ASN:ND2	2.01	0.51
1:A:20:ASN:O	1:A:21:LEU:HD12	2.11	0.51
2:B:450:LYS:HE2	5:B:916:HOH:O	2.11	0.50
2:B:526:GLN:HA	2:B:614:ASN:HD22	1.77	0.50
4:A:301:MRD:H3C1	5:A:454:HOH:O	2.10	0.50
2:B:585:ASP:O	2:B:589:ASN:HB2	2.12	0.50
2:B:392:TYR:HB3	2:B:492:LEU:HB3	1.92	0.50
1:A:93:GLU:O	1:A:96:GLU:HG2	2.11	0.50
4:A:303:MRD:H5C3	2:B:368:ILE:HG23	1.93	0.49
3:C:3:SIA:O1B	3:C:3:SIA:H6	2.13	0.49
2:B:446:ASN:N	5:B:965:HOH:O	2.45	0.49
1:A:165:PHE:HB2	2:B:368:ILE:HD13	1.94	0.49
2:B:427:ILE:HA	2:B:457:GLY:HA3	1.95	0.49
1:A:64:THR:HG21	2:B:287:THR:HG21	1.94	0.49
2:B:458:LEU:CD2	2:B:526:GLN:HB3	2.43	0.49
2:B:336:LYS:HZ2	2:B:338:GLU:HG3	1.78	0.48
1:A:115:MET:HA	1:A:149:LYS:O	2.14	0.48
1:A:108:PRO:CG	1:A:111:LYS:HB2	2.41	0.48
2:B:468:ILE:HD13	2:B:501:ILE:HG13	1.95	0.48
2:B:243:SER:OG	2:B:291:LYS:HE3	2.13	0.48
1:A:40:ASN:H	1:A:40:ASN:ND2	2.11	0.48
2:B:251:ASN:O	2:B:285:THR:HA	2.14	0.48
2:B:386:ILE:HG13	2:B:386:ILE:O	2.14	0.48
2:B:460:ARG:NH2	3:C:1:BGC:H3	2.28	0.48
2:B:550:SER:HA	5:B:845:HOH:O	2.13	0.48
2:B:378:ASN:HD22	2:B:378:ASN:N	2.05	0.47
2:B:444:LEU:C	2:B:444:LEU:HD23	2.39	0.47
2:B:458:LEU:HD22	2:B:526:GLN:HB3	1.97	0.47
2:B:469:VAL:HA	2:B:506:ASP:H	1.80	0.47
1:A:39:GLN:NE2	5:A:413:HOH:O	2.47	0.47
2:B:373:ILE:CD1	2:B:452:ILE:HG21	2.45	0.46
2:B:580:ILE:O	2:B:581:ASN:HB2	2.15	0.46
2:B:555:ARG:NH1	2:B:589:ASN:O	2.47	0.46
2:B:545:ASN:C	2:B:545:ASN:HD22	2.24	0.46
1:A:61:ILE:O	1:A:62:ARG:C	2.58	0.46
1:A:23:ASP:OD1	1:A:130:LEU:HA	2.15	0.46
2:B:545:ASN:HD22	2:B:545:ASN:N	2.13	0.46
2:B:206:ILE:HD12	2:B:206:ILE:N	2.30	0.46
1:A:179:TYR:N	5:A:491:HOH:O	2.34	0.46
2:B:465:SER:HA	2:B:511:ILE:HD12	1.98	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:126:ILE:HB	1:A:138:ALA:HB3	1.97	0.45
1:A:113:LEU:HD11	1:A:150:SER:HB2	1.98	0.45
2:B:497:GLU:HB3	2:B:501:ILE:HG21	1.98	0.45
2:B:526:GLN:HA	2:B:614:ASN:ND2	2.30	0.45
1:A:50:ASN:HB2	2:B:339:VAL:CG2	2.46	0.45
2:B:460:ARG:CZ	3:C:1:BGC:C3	2.94	0.45
2:B:294:GLU:HB3	2:B:304:MET:CE	2.46	0.45
2:B:396:THR:O	2:B:399:GLN:HG3	2.17	0.45
2:B:407:PHE:HE2	2:B:479:ILE:HG13	1.82	0.45
2:B:538:LEU:O	2:B:600:THR:HA	2.16	0.45
2:B:567:ARG:HG3	2:B:567:ARG:HH11	1.81	0.45
2:B:512:LEU:HB2	2:B:538:LEU:HD23	1.99	0.44
2:B:336:LYS:NZ	5:B:854:HOH:O	2.49	0.44
1:A:85:GLN:HE22	1:A:92:THR:N	2.15	0.44
4:A:303:MRD:H5C3	4:A:303:MRD:CM	2.47	0.44
2:B:346:GLN:O	2:B:347:ALA:C	2.61	0.44
1:A:55:ILE:CD1	4:A:301:MRD:H5C3	2.44	0.44
1:A:161:LYS:HB3	1:A:161:LYS:HE2	1.72	0.44
4:B:705:MRD:H5C3	5:B:1060:HOH:O	2.18	0.44
1:A:85:GLN:NE2	1:A:92:THR:N	2.61	0.44
1:A:130:LEU:C	1:A:130:LEU:HD13	2.43	0.44
4:A:302:MRD:HMC2	4:A:302:MRD:O4	2.18	0.44
2:B:430:GLU:HG3	5:B:977:HOH:O	2.18	0.43
4:B:704:MRD:O2	4:B:704:MRD:H5C3	2.18	0.43
4:A:301:MRD:H1C2	4:A:301:MRD:O4	2.17	0.43
4:A:303:MRD:C5	2:B:368:ILE:HG23	2.49	0.43
2:B:418:THR:HB	2:B:466:TYR:HB2	1.98	0.43
2:B:536:TYR:CZ	2:B:617:PHE:HB2	2.53	0.43
1:A:17:ASN:O	1:A:134:VAL:N	2.45	0.43
1:A:60:ARG:O	1:A:60:ARG:HG3	2.19	0.43
1:A:62:ARG:N	5:A:427:HOH:O	2.52	0.43
2:B:373:ILE:HD12	2:B:452:ILE:HG21	2.01	0.43
2:B:439:MET:C	2:B:439:MET:SD	3.02	0.43
2:B:515:LEU:HD11	2:B:538:LEU:CD2	2.49	0.43
1:A:15:SER:N	1:A:18:ASN:HD22	2.16	0.42
1:A:26:TYR:CD1	1:A:181:LYS:HA	2.54	0.42
2:B:517:THR:O	2:B:621:ARG:HD3	2.19	0.42
2:B:551:ILE:CG2	2:B:622:ASN:HB2	2.50	0.42
2:B:424:PRO:HB2	2:B:425:PRO:HD2	2.01	0.42
1:A:40:ASN:HD21	1:A:166:THR:H	1.67	0.42
2:B:545:ASN:ND2	2:B:545:ASN:C	2.77	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:226:ASN:HD21	2:B:355:ASP:N	2.13	0.42
2:B:264:ILE:CG2	2:B:265:PHE:N	2.82	0.42
2:B:204:GLN:N	5:B:1032:HOH:O	2.52	0.42
2:B:237:GLU:O	2:B:240:THR:HB	2.20	0.42
2:B:264:ILE:HD11	4:B:704:MRD:H1C3	2.01	0.42
2:B:336:LYS:CE	2:B:362:THR:HG21	2.48	0.42
2:B:406:LEU:HD13	2:B:478:TYR:CE1	2.54	0.41
2:B:558:ASN:N	5:B:820:HOH:O	2.17	0.41
5:A:421:HOH:O	2:B:436:ASN:HB2	2.20	0.41
2:B:397:SER:HB2	5:B:986:HOH:O	2.20	0.41
1:A:30:ARG:HG3	1:A:125:GLU:OE2	2.21	0.41
2:B:556:THR:HA	2:B:616:ILE:O	2.20	0.41
2:B:555:ARG:HD2	5:B:904:HOH:O	2.19	0.41
1:A:27:VAL:O	1:A:27:VAL:HG13	2.20	0.41
1:A:62:ARG:C	1:A:62:ARG:HE	2.29	0.41
2:B:469:VAL:HG12	2:B:506:ASP:HA	2.02	0.41
2:B:273:ARG:HH11	2:B:273:ARG:HD3	1.72	0.41
2:B:436:ASN:ND2	5:B:1052:HOH:O	2.53	0.41
2:B:527:SER:HA	2:B:528:PRO:HD3	1.92	0.41
1:A:114:TYR:O	1:A:150:SER:HA	2.20	0.41
1:A:154:ASN:OD1	1:A:157:GLU:HG3	2.20	0.41
2:B:316:ILE:HD12	2:B:318:TYR:HB2	2.03	0.41
2:B:500:ASN:C	2:B:502:SER:N	2.77	0.41
2:B:560:GLN:NE2	2:B:587:LEU:HB2	2.36	0.41
1:A:16:ILE:HD13	1:A:16:ILE:O	2.21	0.41
2:B:328:GLU:HG2	5:B:890:HOH:O	2.21	0.41
2:B:540:ILE:CG2	2:B:597:ASN:HA	2.51	0.41
2:B:559:ASN:HB3	2:B:609:GLU:O	2.21	0.41
1:A:38:ARG:HB3	1:A:165:PHE:CZ	2.57	0.40
1:A:128:LYS:HD3	5:A:442:HOH:O	2.20	0.40
2:B:435:SER:HA	2:B:474:ASP:O	2.19	0.40
2:B:545:ASN:HD21	2:B:548:ALA:HB2	1.80	0.40
2:B:545:ASN:O	2:B:545:ASN:CG	2.64	0.40
2:B:496:ARG:NH1	2:B:496:ARG:HG2	2.35	0.40
2:B:447:ASP:CB	2:B:449:ILE:HG13	2.51	0.40
2:B:534:ILE:O	2:B:534:ILE:HG23	2.21	0.40
2:B:296:ALA:HA	2:B:297:PRO:HD3	1.98	0.40
2:B:244:ILE:HD12	2:B:306:VAL:HG23	2.02	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	168/224 (75%)	152 (90%)	14 (8%)	2 (1%)	10	23
2	B	418/420 (100%)	395 (94%)	20 (5%)	3 (1%)	18	38
All	All	586/644 (91%)	547 (93%)	34 (6%)	5 (1%)	14	30

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	546	ASN
2	B	463	ALA
1	A	155	SER
1	A	131	GLN
2	B	547	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	153/207 (74%)	152 (99%)	1 (1%)	76	89
2	B	385/385 (100%)	375 (97%)	10 (3%)	40	68
All	All	538/592 (91%)	527 (98%)	11 (2%)	48	74

All (11) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	16	ILE

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Mol	Chain	Res	Type
2	B	237	GLU
2	B	249	LEU
2	B	316	ILE
2	B	322	ASN
2	B	349	VAL
2	B	373	ILE
2	B	378	ASN
2	B	408	LYS
2	B	545	ASN
2	B	553	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (20) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	39	GLN
1	A	40	ASN
1	A	41	GLN
1	A	82	ASN
1	A	85	GLN
2	B	226	ASN
2	B	275	ASN
2	B	361	GLN
2	B	378	ASN
2	B	393	ASN
2	B	399	GLN
2	B	401	ASN
2	B	483	GLN
2	B	491	GLN
2	B	545	ASN
2	B	546	ASN
2	B	558	ASN
2	B	588	ASN
2	B	612	ASN
2	B	614	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

3 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	BGC	C	1	3	12,12,12	0.39	0	17,17,17	0.68	0
3	GAL	C	2	3	11,11,12	0.70	0	15,15,17	0.64	0
3	SIA	C	3	3	20,20,21	0.77	0	21,28,31	0.88	1 (4%)

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	BGC	C	1	3	-	0/2/22/22	0/1/1/1
3	GAL	C	2	3	-	0/2/19/22	0/1/1/1
3	SIA	C	3	3	-	2/18/34/38	0/1/1/1

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	3	SIA	O1B-C1-C2	2.20	118.43	112.71

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	C	3	SIA	C11-C10-N5-C5

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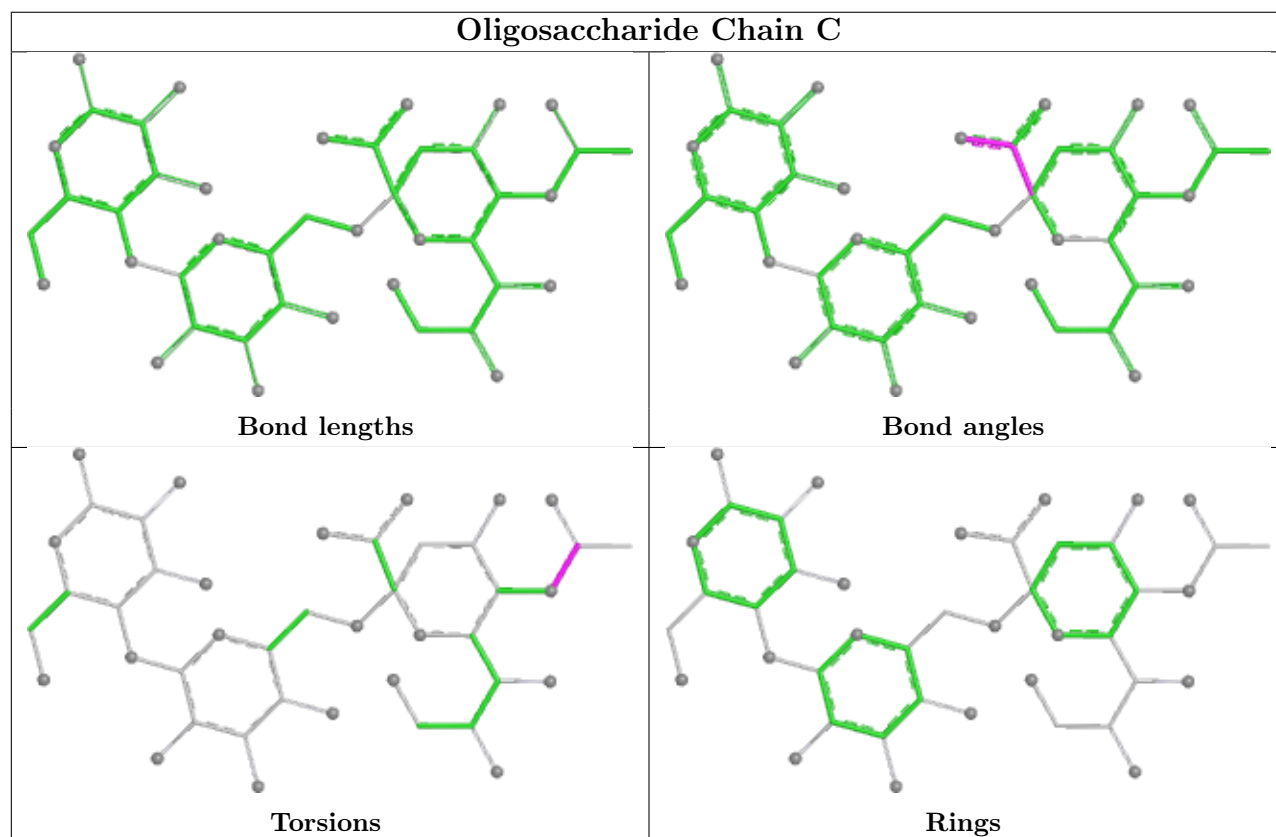
Mol	Chain	Res	Type	Atoms
3	C	3	SIA	O10-C10-N5-C5

There are no ring outliers.

3 monomers are involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	C	2	GAL	1	0
3	C	1	BGC	5	0
3	C	3	SIA	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](#)

5 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	MRD	A	302	-	7,7,7	0.25	0	9,10,10	0.28	0
4	MRD	B	705	-	7,7,7	0.34	0	9,10,10	0.30	0
4	MRD	A	301	-	7,7,7	0.37	0	9,10,10	0.31	0
4	MRD	A	303	-	7,7,7	0.25	0	9,10,10	0.24	0
4	MRD	B	704	-	7,7,7	0.37	0	9,10,10	0.31	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	MRD	A	302	-	-	1/5/5/5	-
4	MRD	B	705	-	-	0/5/5/5	-
4	MRD	A	301	-	-	0/5/5/5	-
4	MRD	A	303	-	-	2/5/5/5	-
4	MRD	B	704	-	-	1/5/5/5	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (4) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	303	MRD	C2-C3-C4-O4
4	A	303	MRD	C2-C3-C4-C5
4	A	302	MRD	O2-C2-C3-C4
4	B	704	MRD	C2-C3-C4-C5

There are no ring outliers.

5 monomers are involved in 20 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	302	MRD	3	0
4	B	705	MRD	2	0
4	A	301	MRD	7	0

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Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	303	MRD	6	0
4	B	704	MRD	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	170/224 (75%)	-0.16	4 (2%) 59 54	28, 48, 74, 86	0
2	B	420/420 (100%)	-0.06	7 (1%) 69 64	29, 51, 71, 86	0
All	All	590/644 (91%)	-0.09	11 (1%) 66 61	28, 50, 72, 86	0

All (11) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	516	ASN	4.2
2	B	547	ILE	3.9
2	B	572	ILE	3.7
1	A	16	ILE	3.2
2	B	546	ASN	2.5
1	A	62	ARG	2.5
1	A	61	ILE	2.3
1	A	184	ASP	2.3
2	B	545	ASN	2.2
2	B	258	ASN	2.2
2	B	515	LEU	2.1

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

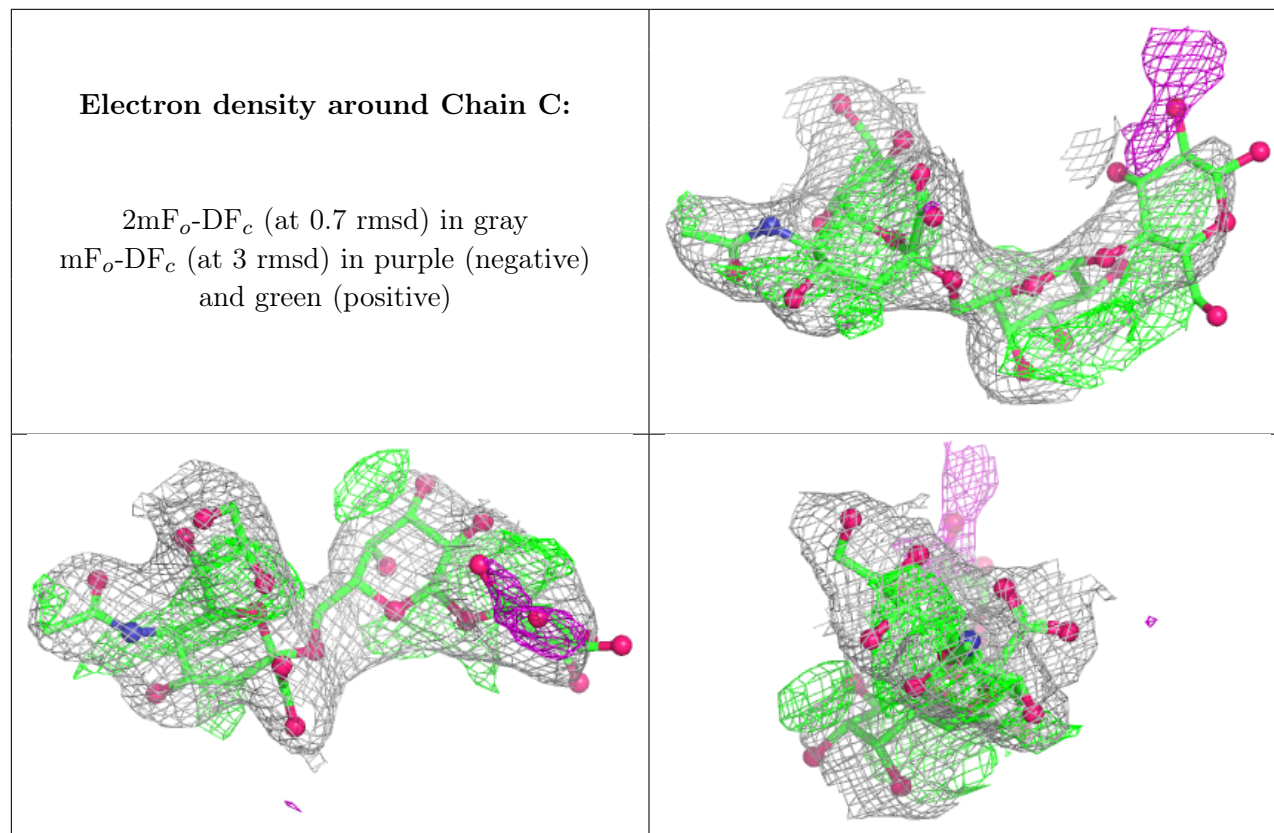
There are no non-standard protein/DNA/RNA residues in this entry.

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($\text{\AA}^2$)	Q<0.9
3	BGC	C	1	12/12	0.60	0.34	78,83,84,84	12
3	GAL	C	2	11/12	0.71	0.19	57,72,76,76	11
3	SIA	C	3	20/21	0.89	0.15	43,48,53,53	20

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
4	MRD	A	302	8/8	0.83	0.38	64,66,66,67	8
4	MRD	B	705	8/8	0.84	0.32	61,62,64,64	8
4	MRD	B	704	8/8	0.85	0.20	55,56,57,59	8
4	MRD	A	303	8/8	0.90	0.27	67,68,68,68	8
4	MRD	A	301	8/8	0.94	0.14	60,61,62,62	8

6.5 Other polymers ⓘ

There are no such residues in this entry.