



# Full wwPDB X-ray Structure Validation Report ⓘ

Aug 9, 2020 – 05:57 AM BST

PDB ID : 4EN6  
Title : Crystal structure of HA70 (HA3) subcomponent of Clostridium botulinum type C progenitor toxin in complex with alpha 2-3-sialyllactose  
Authors : Yamashita, S.; Yoshida, H.; Tonozuka, T.; Nishikawa, A.; Kamitori, S.  
Deposited on : 2012-04-12  
Resolution : 2.56 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](mailto: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 following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467  
Mogul : 1.8.5 (274361), CSD as541be (2020)  
Xtriage (Phenix) : 1.13  
EDS : 2.13.1  
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)  
Refmac : 5.8.0158  
CCP4 : 7.0.044 (Gargrove)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.13.1

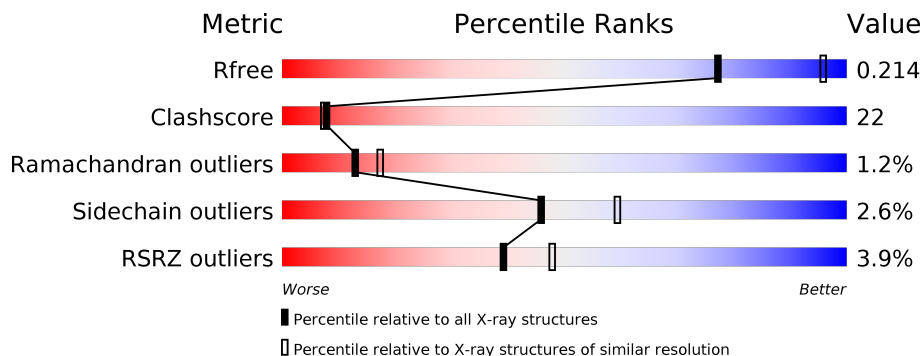
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.56 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))
$R_{free}$	130704	1279 (2.58-2.54)
Clashscore	141614	1327 (2.58-2.54)
Ramachandran outliers	138981	1312 (2.58-2.54)
Sidechain outliers	138945	1312 (2.58-2.54)
RSRZ outliers	127900	1269 (2.58-2.54)

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 on the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 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	<p>4% 47% 26% 24%</p>
2	B	420	<p>4% 70% 28%</p>
3	C	3	<p>67% 33%</p>

## 2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 5203 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
			Total	C	N	O	S			
1	A	170	1382	882	226	271	3	0	0	0

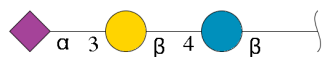
There are 21 discrepancies between the modelled and reference sequences:

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.

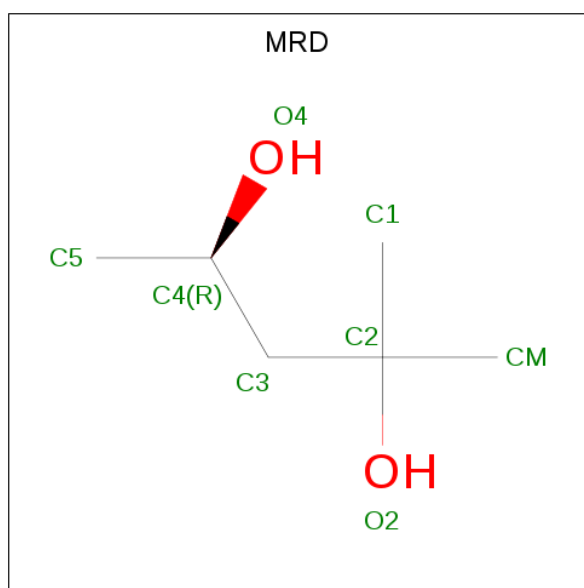
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	B	420	3339	2109	553	672	5	0	0	0

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



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

- Molecule 4 is (4R)-2-METHYLPENTANE-2,4-DIOL (three-letter code: MRD) (formula: C<sub>6</sub>H<sub>14</sub>O<sub>2</sub>).



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		

*Continued on next page...*

*Continued from previous page...*

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

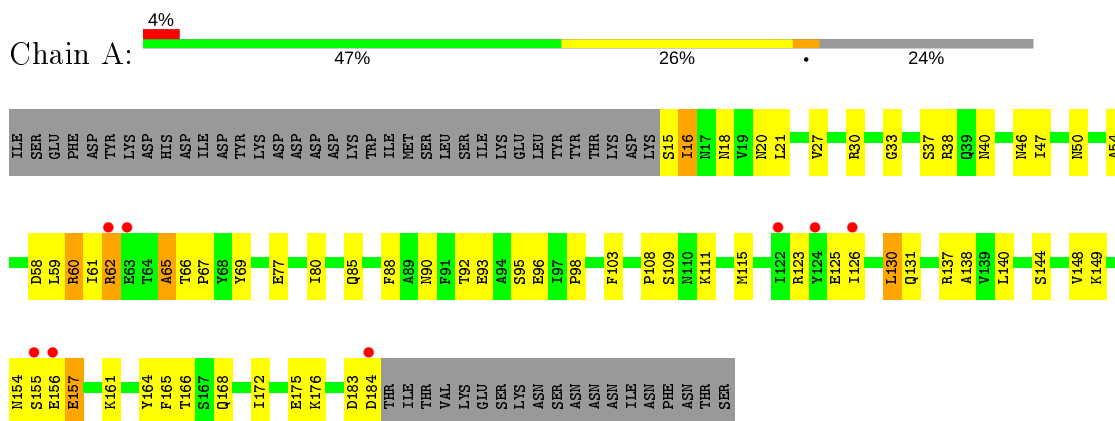
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	106	Total	O	0	0
			106	106		
5	B	293	Total	O	0	0
			293	293		

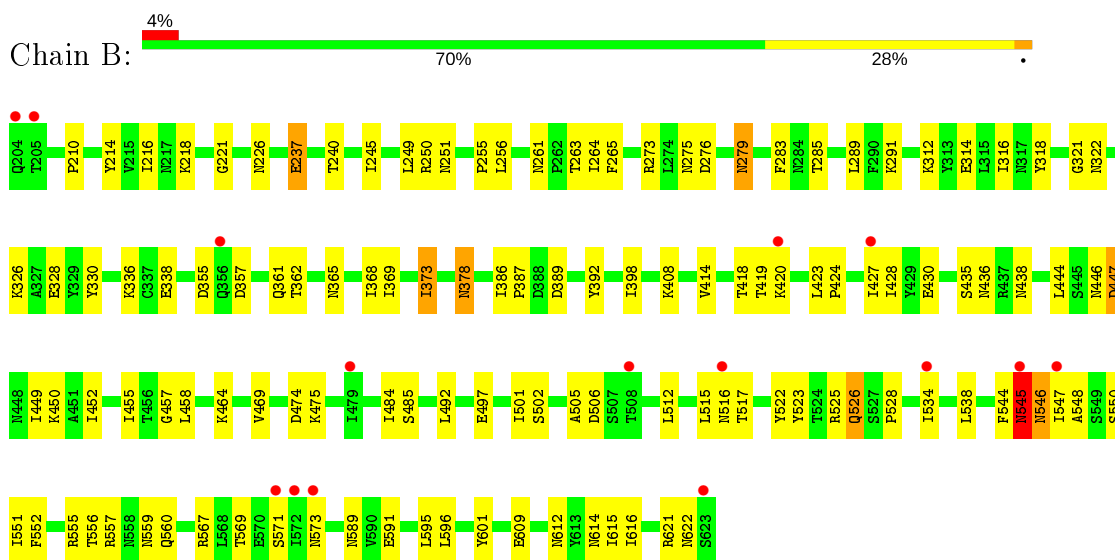
### 3 Residue-property plots [i](#)

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: Hemagglutinin components HA-22/23/53



- Molecule 2: Hemagglutinin components HA-22/23/53



- Molecule 3: N-acetyl-alpha-neuraminic acid-(2-3)-beta-D-galactopyranose-(1-4)-beta-D-glucopyranose





## 4 Data and refinement statistics

Property	Value	Source
Space group	P 63	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	175.60 Å 175.60 Å 80.78 Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	25.74 – 2.56 25.74 – 2.56	Depositor EDS
% Data completeness (in resolution range)	97.3 (25.74-2.56) 97.4 (25.74-2.56)	Depositor EDS
$R_{merge}$	0.09	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	3.95 (at 2.57 Å)	Xtrriage
Refinement program	CNS 1.1	Depositor
R, $R_{free}$	0.188 , 0.219 0.182 , 0.214	Depositor DCC
$R_{free}$ test set	4499 reflections (9.81%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	48.4	Xtrriage
Anisotropy	0.082	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.32 , 51.6	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.33$	Xtrriage
Estimated twinning fraction	0.028 for h,-h-k,-l	Xtrriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	5203	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	52.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 10.83% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>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

### 5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: MRD, SIA, BGC, GAL

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	0.37	0/1411	0.66	0/1914
2	B	0.34	0/3405	0.67	0/4642
All	All	0.35	0/4816	0.67	0/6556

There are no bond length outliers.

There are no bond angle outliers.

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	63	0
2	B	3339	0	3275	138	0
3	C	43	0	37	3	0
4	A	24	0	42	12	0
4	B	16	0	28	4	0
5	A	106	0	0	13	0
5	B	293	0	0	29	0
All	All	5203	0	4718	206	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 22.

All (206) 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	1.57	0.84
2:B:264:ILE:HD11	4:B:704:MRD:HMC3	1.60	0.84
2:B:612:ASN:ND2	2:B:614:ASN:H	1.78	0.81
2:B:435:SER:HB3	2:B:475:LYS:HG2	1.61	0.81
2:B:273:ARG:NH1	2:B:276:ASP:OD2	2.14	0.79
2:B:264:ILE:HG23	2:B:265:PHE:H	1.49	0.77
2:B:226:ASN:HD21	2:B:355:ASP:H	1.32	0.77
2:B:221:GLY:O	2:B:362:THR:HB	1.84	0.77
2:B:545:ASN:HD21	2:B:548:ALA:HB3	1.48	0.77
2:B:612:ASN:HD22	2:B:614:ASN:H	1.32	0.74
1:A:37:SER:O	1:A:38:ARG:HG2	1.85	0.74
2:B:216:ILE:HD12	2:B:369:ILE:HG12	1.69	0.74
2:B:336:LYS:HE2	2:B:362:THR:HG23	1.72	0.72
2:B:373:ILE:HD12	2:B:452:ILE:HD13	1.71	0.72
2:B:484:ILE:HB	5:B:876:HOH:O	1.90	0.72
2:B:545:ASN:HD21	2:B:548:ALA:CB	2.03	0.71
2:B:567:ARG:HD3	2:B:601:TYR:CE1	2.26	0.71
4:A:303:MRD:HMC1	4:A:303:MRD:H5C3	1.71	0.70
2:B:336:LYS:HE2	2:B:362:THR:CG2	2.22	0.69
1:A:85:GLN:NE2	1:A:92:THR:H	1.91	0.69
1:A:108:PRO:HG2	1:A:111:LYS:HB2	1.74	0.68
2:B:378:ASN:H	2:B:378:ASN:HD22	1.42	0.68
2:B:261:ASN:OD1	2:B:264:ILE:HG22	1.94	0.67
1:A:176:LYS:HE3	5:A:476:HOH:O	1.95	0.67
2:B:545:ASN:ND2	2:B:548:ALA:HB3	2.11	0.66
2:B:555:ARG:HD3	5:B:884:HOH:O	1.95	0.65
2:B:214:TYR:HB2	2:B:316:ILE:HD11	1.78	0.64
2:B:314:GLU:OE2	2:B:326:LYS:HE3	1.97	0.64
2:B:256:LEU:HB2	2:B:330:TYR:HB2	1.80	0.64
2:B:264:ILE:HG21	5:B:960:HOH:O	1.96	0.64
2:B:545:ASN:HD22	2:B:545:ASN:N	1.95	0.64
1:A:183:ASP:CG	1:A:184:ASP:H	2.00	0.64
1:A:40:ASN:HD22	1:A:40:ASN:H	1.46	0.63
1:A:47:ILE:HD13	2:B:362:THR:HG22	1.79	0.63
4:A:301:MRD:H1C1	5:A:471:HOH:O	1.98	0.63
2:B:336:LYS:HZ2	2:B:338:GLU:HG3	1.64	0.63
2:B:505:ALA:HB3	5:B:897:HOH:O	1.99	0.62
2:B:216:ILE:CD1	2:B:369:ILE:HG23	2.29	0.62
1:A:59:LEU:O	1:A:61:ILE:N	2.33	0.61
4:A:301:MRD:H3C1	5:A:471:HOH:O	1.99	0.61

*Continued on next page...*

*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:B:705:MRD:H5C3	5:B:1027:HOH:O	1.99	0.61
1:A:161:LYS:HG2	5:A:493:HOH:O	1.99	0.61
1:A:184:ASP:HA	5:A:430:HOH:O	2.00	0.61
2:B:250:ARG:HA	2:B:250:ARG:NE	2.16	0.61
2:B:264:ILE:HG23	2:B:265:PHE:N	2.14	0.61
1:A:77:GLU:HB2	5:A:470:HOH:O	2.01	0.61
4:A:303:MRD:C5	2:B:368:ILE:HG23	2.31	0.60
2:B:516:ASN:HA	2:B:621:ARG:NH1	2.15	0.60
1:A:168:GLN:HB2	5:A:460:HOH:O	2.02	0.59
2:B:552:PHE:CD2	2:B:595:LEU:HD12	2.37	0.59
2:B:595:LEU:HB3	2:B:601:TYR:CE2	2.36	0.59
2:B:428:ILE:HG23	2:B:455:ILE:HB	1.84	0.59
2:B:336:LYS:NZ	2:B:338:GLU:HG3	2.17	0.59
2:B:447:ASP:HB2	2:B:449:ILE:HG13	1.85	0.59
4:A:302:MRD:H5C3	2:B:289:LEU:HD22	1.84	0.59
2:B:551:ILE:HB	2:B:622:ASN:HB2	1.84	0.59
1:A:165:PHE:HB2	2:B:368:ILE:HD13	1.84	0.59
2:B:361:GLN:NE2	5:B:1000:HOH:O	2.34	0.59
2:B:528:PRO:HG2	2:B:534:ILE:HB	1.84	0.58
1:A:60:ARG:O	1:A:60:ARG:HG3	2.03	0.58
2:B:423:LEU:HD13	2:B:464:LYS:HA	1.85	0.58
2:B:591:GLU:HG2	5:B:1008:HOH:O	2.03	0.58
1:A:65:ALA:HA	5:A:497:HOH:O	2.05	0.57
1:A:98:PRO:HA	5:A:473:HOH:O	2.03	0.57
1:A:93:GLU:O	1:A:96:GLU:HG2	2.05	0.57
1:A:66:THR:HG23	5:A:443:HOH:O	2.04	0.57
1:A:16:ILE:HD13	1:A:16:ILE:O	2.05	0.57
1:A:60:ARG:NH2	5:A:468:HOH:O	2.37	0.57
4:A:303:MRD:H5C3	2:B:368:ILE:HG23	1.87	0.57
3:C:3:SIA:O1B	3:C:3:SIA:H6	2.04	0.56
1:A:90:ASN:OD1	1:A:93:GLU:HG3	2.06	0.56
2:B:512:LEU:HB2	2:B:538:LEU:HD23	1.89	0.55
2:B:559:ASN:HB3	2:B:609:GLU:O	2.07	0.55
2:B:218:LYS:HE2	5:B:958:HOH:O	2.06	0.55
2:B:469:VAL:HG12	2:B:506:ASP:HA	1.89	0.55
2:B:515:LEU:CD1	2:B:538:LEU:HD22	2.37	0.55
1:A:40:ASN:ND2	1:A:40:ASN:H	2.04	0.55
2:B:545:ASN:O	2:B:547:ILE:N	2.40	0.55
2:B:444:LEU:O	2:B:444:LEU:HD23	2.07	0.55
1:A:85:GLN:HE22	1:A:92:THR:H	1.55	0.55
2:B:447:ASP:CB	2:B:449:ILE:HG13	2.37	0.54

*Continued on next page...*

*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:484:ILE:O	2:B:485:SER:HB3	2.07	0.54
2:B:515:LEU:HD11	2:B:538:LEU:HD22	1.89	0.54
4:A:303:MRD:H5C2	5:B:877:HOH:O	2.07	0.54
1:A:88:PHE:HB3	5:A:405:HOH:O	2.07	0.54
2:B:612:ASN:HD22	2:B:614:ASN:N	2.04	0.54
2:B:450:LYS:HE2	5:B:892:HOH:O	2.08	0.54
2:B:528:PRO:CG	2:B:534:ILE:HB	2.38	0.54
2:B:550:SER:HA	5:B:1070:HOH:O	2.07	0.53
2:B:261:ASN:CG	2:B:264:ILE:HG22	2.29	0.53
2:B:328:GLU:HG2	5:B:1007:HOH:O	2.09	0.53
1:A:69:TYR:CE1	1:A:140:LEU:HD13	2.44	0.52
2:B:557:ARG:HG2	2:B:557:ARG:HH11	1.74	0.52
1:A:108:PRO:CG	1:A:111:LYS:HB2	2.38	0.52
2:B:485:SER:N	5:B:876:HOH:O	2.36	0.52
2:B:512:LEU:HG	2:B:523:TYR:CE2	2.45	0.52
2:B:256:LEU:HD11	5:B:1088:HOH:O	2.09	0.52
2:B:458:LEU:HD22	2:B:526:GLN:HB3	1.92	0.52
2:B:447:ASP:HB3	5:B:1012:HOH:O	2.08	0.51
2:B:264:ILE:HB	5:B:930:HOH:O	2.10	0.51
2:B:336:LYS:CD	2:B:362:THR:HG21	2.40	0.51
2:B:420:LYS:HD3	5:B:963:HOH:O	2.10	0.51
2:B:555:ARG:NH1	2:B:589:ASN:O	2.44	0.51
1:A:37:SER:C	1:A:38:ARG:HG2	2.31	0.51
1:A:59:LEU:HD11	4:A:302:MRD:CM	2.41	0.50
2:B:545:ASN:O	2:B:546:ASN:C	2.50	0.50
1:A:50:ASN:HA	1:A:109:SER:O	2.12	0.50
2:B:237:GLU:O	2:B:240:THR:HB	2.12	0.50
2:B:547:ILE:HG13	2:B:548:ALA:N	2.27	0.50
1:A:154:ASN:OD1	1:A:157:GLU:HG3	2.11	0.50
2:B:512:LEU:CD2	2:B:538:LEU:HD21	2.42	0.49
2:B:387:PRO:HA	5:B:1035:HOH:O	2.13	0.49
2:B:386:ILE:O	2:B:386:ILE:HG13	2.12	0.49
2:B:392:TYR:HB3	2:B:492:LEU:HB3	1.95	0.49
1:A:154:ASN:O	1:A:156:GLU:N	2.46	0.49
2:B:398:ILE:HG12	2:B:484:ILE:HG12	1.94	0.49
2:B:515:LEU:HD11	2:B:538:LEU:CD2	2.42	0.49
1:A:30:ARG:HB2	1:A:123:ARG:HB3	1.95	0.49
2:B:560:GLN:HB3	2:B:609:GLU:HB3	1.93	0.49
2:B:365:ASN:HB3	5:B:1030:HOH:O	2.12	0.48
4:A:301:MRD:O4	4:A:301:MRD:H1C2	2.12	0.48
2:B:435:SER:HA	2:B:474:ASP:O	2.13	0.48

*Continued on next page...*

*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:20:ASN:O	1:A:21:LEU:HD12	2.14	0.48
1:A:130:LEU:CD1	1:A:131:GLN:HG2	2.43	0.48
2:B:318:TYR:OH	2:B:321:GLY:HA2	2.13	0.48
1:A:85:GLN:HE22	1:A:92:THR:CB	2.26	0.48
1:A:15:SER:HA	5:A:449:HOH:O	2.14	0.48
2:B:559:ASN:HB2	2:B:615:ILE:HD13	1.96	0.48
1:A:183:ASP:CG	1:A:184:ASP:N	2.65	0.48
2:B:336:LYS:HD2	2:B:362:THR:HG21	1.96	0.48
4:A:302:MRD:HMC2	4:A:302:MRD:O4	2.14	0.47
1:A:46:ASN:ND2	1:A:54:ALA:HB1	2.30	0.47
1:A:16:ILE:CD1	1:A:16:ILE:O	2.63	0.47
2:B:556:THR:HA	2:B:616:ILE:O	2.14	0.47
2:B:567:ARG:NH1	2:B:569:THR:HG22	2.30	0.47
1:A:30:ARG:NH1	1:A:175:GLU:OE2	2.44	0.47
1:A:115:MET:HA	1:A:149:LYS:O	2.15	0.47
2:B:398:ILE:CG1	2:B:484:ILE:HG12	2.46	0.46
2:B:526:GLN:NE2	5:B:1055:HOH:O	2.48	0.46
1:A:103:PHE:CD2	1:A:148:VAL:HG13	2.51	0.46
1:A:40:ASN:HD21	1:A:166:THR:H	1.63	0.46
1:A:130:LEU:O	1:A:131:GLN:HB2	2.16	0.46
2:B:458:LEU:HD13	2:B:525:ARG:HD2	1.97	0.46
1:A:38:ARG:HB3	1:A:165:PHE:CZ	2.51	0.45
1:A:126:ILE:HB	1:A:138:ALA:HB3	1.97	0.45
2:B:444:LEU:C	2:B:444:LEU:HD23	2.36	0.45
2:B:250:ARG:CZ	2:B:250:ARG:HA	2.46	0.45
2:B:512:LEU:HD23	2:B:538:LEU:HD21	1.98	0.45
2:B:427:ILE:HA	2:B:457:GLY:HA3	1.98	0.45
1:A:69:TYR:CD1	1:A:140:LEU:HD13	2.52	0.45
2:B:336:LYS:HE2	2:B:362:THR:HG21	1.98	0.45
2:B:430:GLU:HG2	5:B:948:HOH:O	2.16	0.45
2:B:419:THR:HG21	2:B:424:PRO:HD2	1.98	0.45
1:A:15:SER:N	1:A:18:ASN:ND2	2.64	0.44
1:A:125:GLU:OE1	1:A:137:ARG:NH2	2.47	0.44
2:B:289:LEU:HD12	2:B:289:LEU:N	2.32	0.44
2:B:414:VAL:HG21	2:B:502:SER:CB	2.47	0.44
2:B:378:ASN:ND2	2:B:378:ASN:H	2.12	0.44
2:B:214:TYR:HB2	2:B:316:ILE:CD1	2.47	0.44
1:A:58:ASP:OD2	1:A:59:LEU:N	2.45	0.44
2:B:387:PRO:HG2	2:B:389:ASP:O	2.17	0.44
1:A:156:GLU:OE1	1:A:156:GLU:N	2.51	0.44
2:B:264:ILE:CG2	2:B:265:PHE:H	2.27	0.43

*Continued on next page...*

*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:263:THR:HG22	5:B:895:HOH:O	2.18	0.43
1:A:80:ILE:HD11	1:A:138:ALA:HB1	2.01	0.43
2:B:318:TYR:CZ	2:B:321:GLY:HA2	2.53	0.43
1:A:59:LEU:HD11	4:A:302:MRD:HMC2	1.99	0.43
1:A:60:ARG:HD3	1:A:144:SER:O	2.19	0.43
2:B:418:THR:HG22	5:B:962:HOH:O	2.18	0.43
2:B:251:ASN:O	2:B:285:THR:HA	2.19	0.43
1:A:137:ARG:O	2:B:438:ASN:HA	2.18	0.43
1:A:33:GLY:O	1:A:172:ILE:HG12	2.18	0.43
2:B:264:ILE:HD11	4:B:704:MRD:CM	2.40	0.43
2:B:275:ASN:O	2:B:279:ASN:HA	2.18	0.43
2:B:373:ILE:CD1	2:B:452:ILE:HG21	2.49	0.43
2:B:414:VAL:HG21	2:B:502:SER:HB3	2.01	0.43
2:B:544:PHE:HE1	2:B:596:LEU:HD23	1.84	0.43
1:A:20:ASN:C	1:A:21:LEU:HD12	2.39	0.42
2:B:210:PRO:HA	2:B:526:GLN:HE22	1.84	0.42
2:B:245:ILE:HG13	2:B:291:LYS:HG3	2.01	0.42
2:B:436:ASN:HB3	5:B:994:HOH:O	2.19	0.42
2:B:497:GLU:HB3	2:B:501:ILE:HG21	2.01	0.42
2:B:523:TYR:CE1	3:C:3:SIA:H112	2.54	0.42
2:B:256:LEU:HA	2:B:256:LEU:HD23	1.92	0.42
2:B:312:LYS:HE2	5:B:1007:HOH:O	2.18	0.41
2:B:516:ASN:HA	2:B:621:ARG:HH12	1.84	0.41
2:B:517:THR:O	2:B:621:ARG:HD3	2.21	0.41
2:B:458:LEU:CD2	2:B:526:GLN:HB3	2.50	0.41
2:B:545:ASN:ND2	2:B:545:ASN:N	2.66	0.41
2:B:273:ARG:NH2	5:B:875:HOH:O	2.48	0.41
1:A:62:ARG:HB2	1:A:95:SER:O	2.21	0.41
2:B:256:LEU:HD21	5:B:1088:HOH:O	2.20	0.41
4:B:704:MRD:H5C3	4:B:704:MRD:O2	2.20	0.41
1:A:66:THR:HA	1:A:67:PRO:HD3	1.93	0.41
2:B:226:ASN:HD21	2:B:355:ASP:N	2.08	0.41
1:A:164:TYR:HB2	2:B:368:ILE:HD11	2.02	0.41
1:A:27:VAL:HG13	1:A:27:VAL:O	2.20	0.41
1:A:47:ILE:CD1	2:B:362:THR:HG22	2.46	0.41
3:C:3:SIA:C6	3:C:3:SIA:O1B	2.69	0.41
2:B:525:ARG:C	2:B:614:ASN:HD22	2.23	0.41
2:B:526:GLN:NE2	5:B:1054:HOH:O	2.53	0.41
2:B:255:PRO:HG3	2:B:283:PHE:O	2.20	0.41
2:B:559:ASN:HB2	2:B:615:ILE:CD1	2.51	0.41
2:B:571:SER:C	2:B:573:ASN:H	2.23	0.41

*Continued on next page...*

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:273:ARG:NE	5:B:875:HOH:O	2.44	0.40
2:B:522:TYR:CE1	2:B:557:ARG:HD3	2.56	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	168/224 (75%)	148 (88%)	15 (9%)	5 (3%)	<b>4</b> <b>4</b>
2	B	418/420 (100%)	397 (95%)	19 (4%)	2 (0%)	29 39
All	All	586/644 (91%)	545 (93%)	34 (6%)	7 (1%)	<b>13</b> <b>17</b>

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	60	ARG
1	A	155	SER
1	A	157	GLU
2	B	546	ASN
2	B	545	ASN
1	A	65	ALA
1	A	62	ARG

### 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	153/207 (74%)	151 (99%)	2 (1%)	69	80
2	B	385/385 (100%)	373 (97%)	12 (3%)	40	52
All	All	538/592 (91%)	524 (97%)	14 (3%)	46	59

All (14) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	16	ILE
1	A	130	LEU
2	B	237	GLU
2	B	249	LEU
2	B	279	ASN
2	B	322	ASN
2	B	357	ASP
2	B	373	ILE
2	B	378	ASN
2	B	408	LYS
2	B	446	ASN
2	B	447	ASP
2	B	526	GLN
2	B	545	ASN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (15) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	40	ASN
1	A	41	GLN
1	A	46	ASN
1	A	82	ASN
1	A	85	GLN
2	B	226	ASN
2	B	279	ASN
2	B	378	ASN
2	B	399	GLN
2	B	422	ASN
2	B	545	ASN
2	B	558	ASN
2	B	576	ASN
2	B	612	ASN
2	B	614	ASN



### 5.3.3 RNA [i](#)

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	0.40	0	17,17,17	0.51	0
3	GAL	C	2	3	11,11,12	0.36	0	15,15,17	0.58	0
3	SIA	C	3	3	17,20,21	0.58	0	21,28,31	0.70	0

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	BGC	C	1	3	-	2/2/22/22	0/1/1/1
3	GAL	C	2	3	-	0/2/19/22	0/1/1/1
3	SIA	C	3	3	-	0/14/34/38	0/1/1/1

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (2) torsion outliers are listed below:

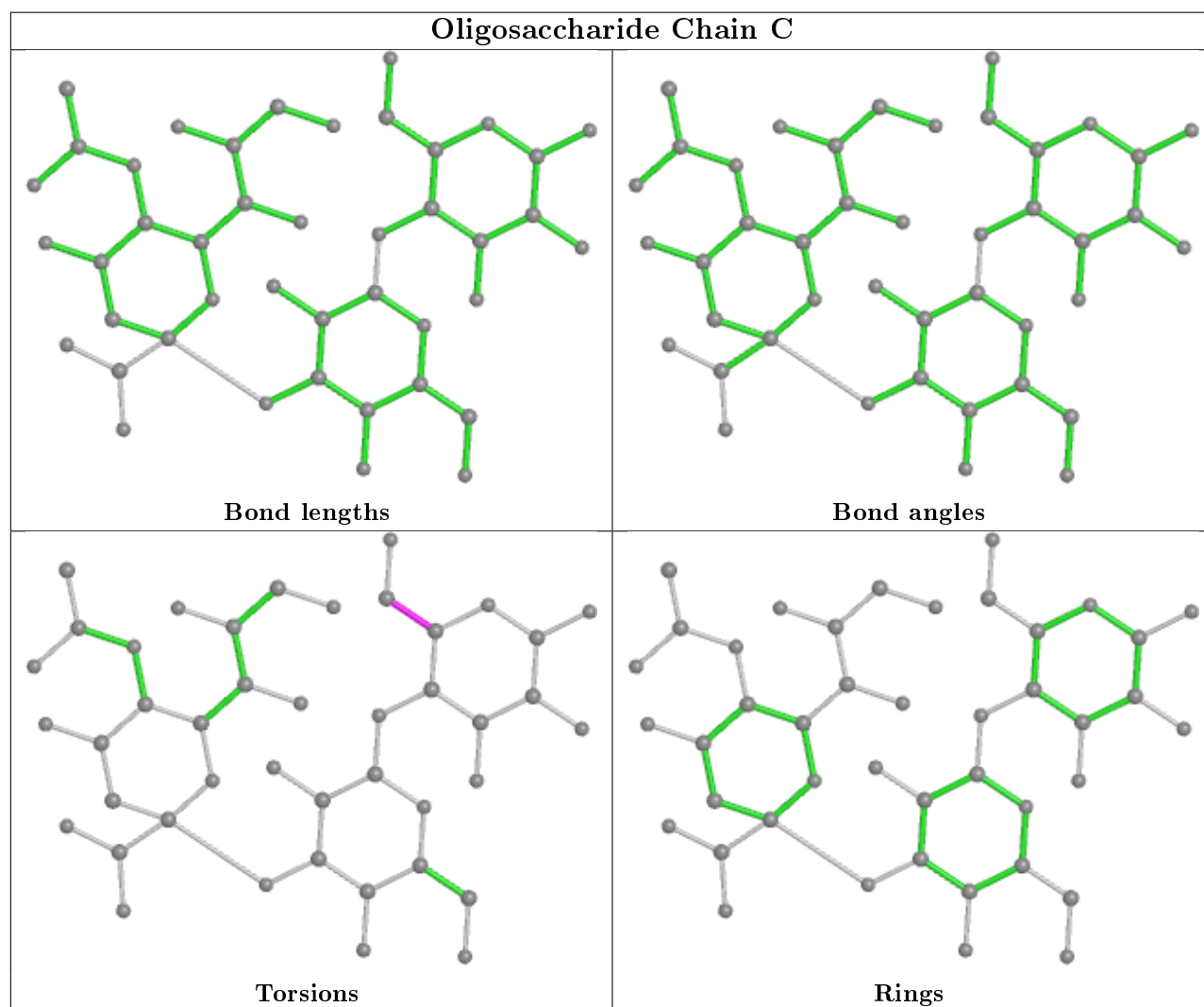
Mol	Chain	Res	Type	Atoms
3	C	1	BGC	O5-C5-C6-O6
3	C	1	BGC	C4-C5-C6-O6

There are no ring outliers.

1 monomer is involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	C	3	SIA	3	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	301	-	7,7,7	0.25	0	9,10,10	0.34	0
4	MRD	A	302	-	7,7,7	0.24	0	9,10,10	0.25	0
4	MRD	A	303	-	7,7,7	0.24	0	9,10,10	0.28	0
4	MRD	B	704	-	7,7,7	0.29	0	9,10,10	0.29	0
4	MRD	B	705	-	7,7,7	0.23	0	9,10,10	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	MRD	A	301	-	-	0/5/5/5	-
4	MRD	A	302	-	-	0/5/5/5	-
4	MRD	A	303	-	-	1/5/5/5	-
4	MRD	B	704	-	-	1/5/5/5	-
4	MRD	B	705	-	-	1/5/5/5	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (3) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	303	MRD	C2-C3-C4-O4
4	B	704	MRD	C2-C3-C4-O4
4	B	705	MRD	C2-C3-C4-C5

There are no ring outliers.

5 monomers are involved in 16 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	301	MRD	3	0

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	302	MRD	4	0
4	A	303	MRD	5	0
4	B	704	MRD	3	0
4	B	705	MRD	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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	170/224 (75%)	0.01	8 (4%) 31 40	29, 50, 80, 90	0
2	B	420/420 (100%)	-0.06	15 (3%) 42 51	27, 50, 74, 90	0
All	All	590/644 (91%)	-0.04	23 (3%) 39 48	27, 50, 76, 90	0

All (23) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	547	ILE	5.4
2	B	572	ILE	4.6
2	B	204	GLN	4.2
1	A	62	ARG	3.6
2	B	573	ASN	3.5
1	A	63	GLU	3.5
2	B	205	THR	3.4
2	B	623	SER	3.3
1	A	124	TYR	3.1
2	B	420	LYS	2.9
2	B	571	SER	2.8
1	A	122	ILE	2.8
2	B	516	ASN	2.7
2	B	479	ILE	2.6
1	A	156	GLU	2.5
1	A	126	ILE	2.3
2	B	427	ILE	2.2
2	B	508	THR	2.2
2	B	356	GLN	2.2
1	A	155	SER	2.2
2	B	545	ASN	2.1
1	A	184	ASP	2.0
2	B	534	ILE	2.0

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

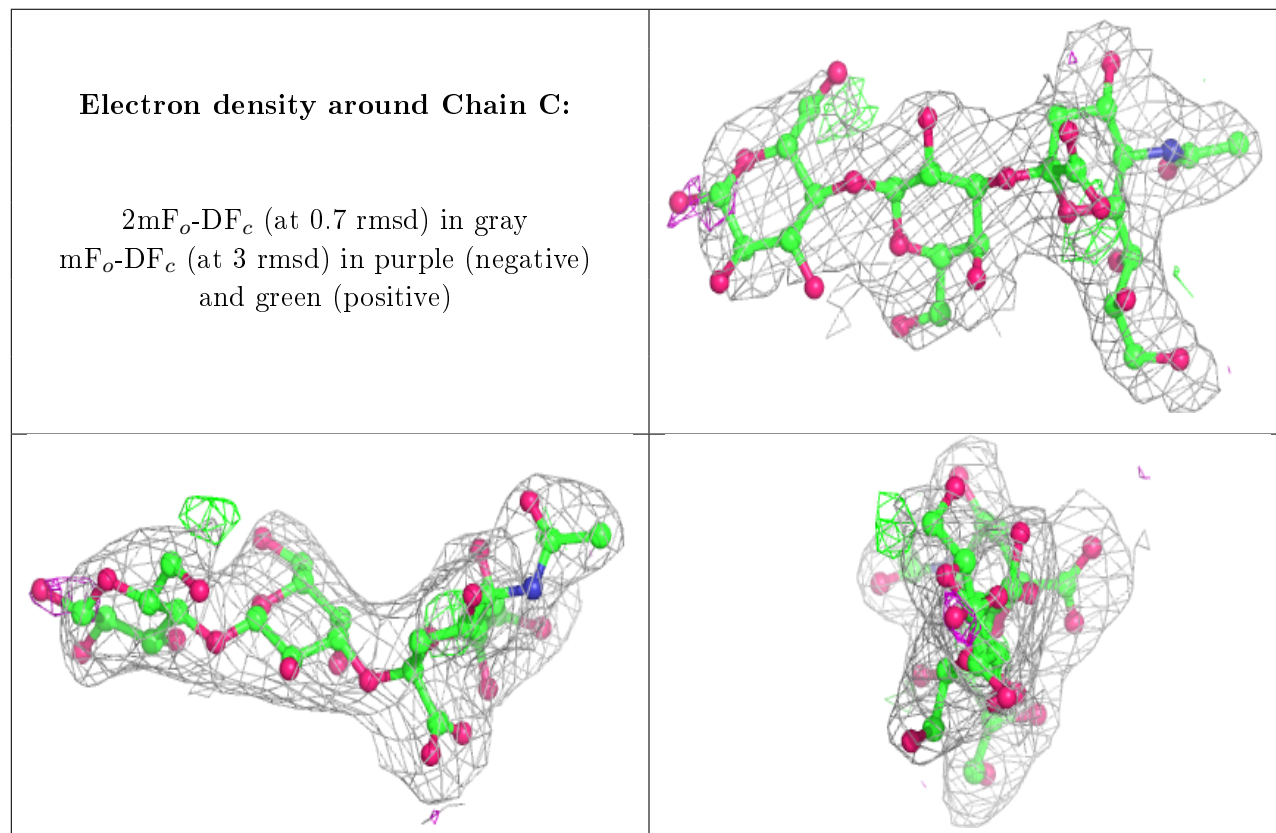
There are no non-standard protein/DNA/RNA residues in this entry.

## 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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
3	BGC	C	1	12/12	0.84	0.46	74,85,87,89	0
3	SIA	C	3	20/21	0.95	0.12	47,51,54,55	0
3	GAL	C	2	11/12	0.96	0.21	57,66,68,69	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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
4	MRD	A	302	8/8	0.82	0.53	71,72,76,77	0
4	MRD	B	705	8/8	0.86	0.47	68,70,71,73	0
4	MRD	B	704	8/8	0.90	0.20	61,63,64,67	0
4	MRD	A	301	8/8	0.91	0.31	63,65,66,67	0
4	MRD	A	303	8/8	0.91	0.34	70,71,73,73	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.