



Full wwPDB X-ray Structure Validation Report ⓘ

Aug 10, 2020 – 01:27 AM BST

PDB ID : 2IY8
Title : Crystal structure of the sialyltransferase PM0188 with CMP-3FNeuAc and lactose
Authors : Kim, D.U.; Cho, H.S.
Deposited on : 2006-07-13
Resolution : 2.50 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 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
buster-report : 1.1.7 (2018)
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

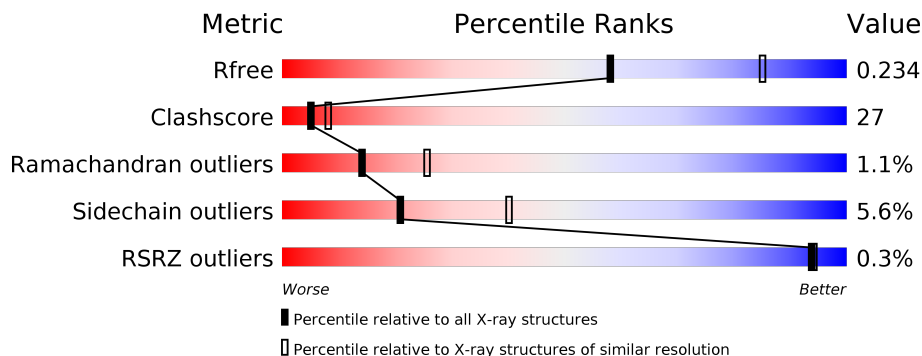
1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 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	4661 (2.50-2.50)
Clashscore	141614	5346 (2.50-2.50)
Ramachandran outliers	138981	5231 (2.50-2.50)
Sidechain outliers	138945	5233 (2.50-2.50)
RSRZ outliers	127900	4559 (2.50-2.50)

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 ≥ 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	389	 55% 37%
2	B	2	 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
3	CSF	A	1415	X	-	-	-

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 3302 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 PROTEIN PM0188.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	Se			
1	A	377	3107	2007	511	583	6	0	0	0

There are 5 discrepancies between the modelled and reference sequences:

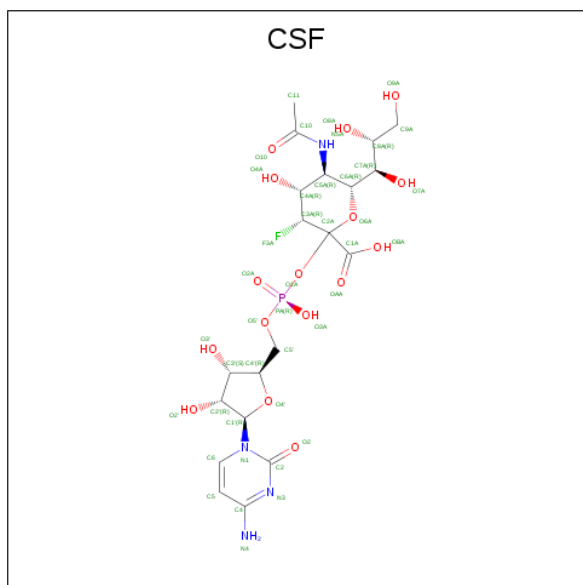
Chain	Residue	Modelled	Actual	Comment	Reference
A	105	ASN	ASP	conflict	UNP Q9CP67
A	135	GLN	ARG	conflict	UNP Q9CP67
A	275	GLU	ASP	conflict	UNP Q9CP67
A	295	GLU	GLY	conflict	UNP Q9CP67
A	411	GLU	GLN	conflict	UNP Q9CP67

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



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace
			Total	C	O			
2	B	2	23	12	11	0	0	0

- Molecule 3 is CYTIDINE-5'-MONOPHOSPHATE-3-FLUORO-N-ACETYL-NEURAMINI C ACID (three-letter code: CSF) (formula: C₂₀H₃₀FN₄O₁₆P).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	
			Total	C	F	N	O			P
3	A	1	42	20	1	4	16	1	0	0

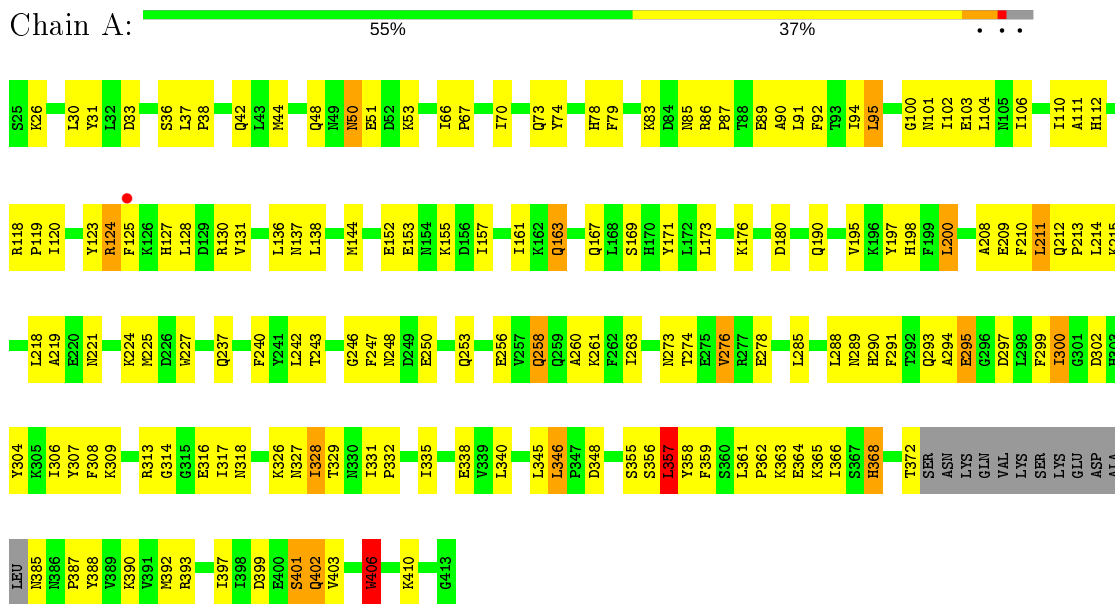
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	130	Total	O	0	0
			130	130		

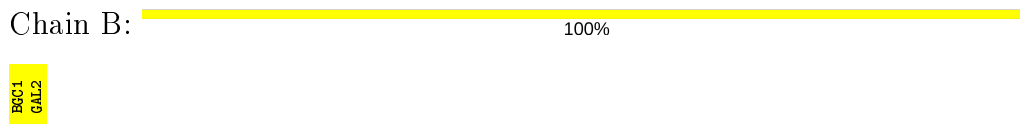
3 Residue-property plots [i](#)

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: PROTEIN PM0188



- Molecule 2: beta-D-galactopyranose-(1-4)-beta-D-glucopyranose



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	64.74Å 65.45Å 105.64Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	15.00 – 2.50 28.15 – 2.50	Depositor EDS
% Data completeness (in resolution range)	91.3 (15.00-2.50) 91.3 (28.15-2.50)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.71 (at 2.51Å)	Xtrriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.235 , 0.244 0.225 , 0.234	Depositor DCC
R_{free} test set	753 reflections (5.11%)	wwPDB-VP
Wilson B-factor (Å ²)	19.5	Xtrriage
Anisotropy	0.341	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 1.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.28$	Xtrriage
Estimated twinning fraction	0.056 for k,h,-l	Xtrriage
F_o, F_c correlation	0.88	EDS
Total number of atoms	3302	wwPDB-VP
Average B, all atoms (Å ²)	17.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.81% 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 [i](#)

5.1 Standard geometry [i](#)

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

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	0.41	0/3177	0.66	1/4296 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	357	LEU	CA-CB-CG	6.42	130.08	115.30

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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	3107	0	3052	169	0
2	B	23	0	21	0	0
3	A	42	0	26	2	0
4	A	130	0	0	23	2
All	All	3302	0	3099	169	2

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 27.

All (169) 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:273:ASN:HB3	4:A:2089:HOH:O	1.29	1.33
1:A:209:GLU:CD	4:A:2068:HOH:O	1.78	1.17
1:A:209:GLU:OE2	4:A:2068:HOH:O	1.57	1.16
1:A:209:GLU:OE1	4:A:2068:HOH:O	1.68	1.05
1:A:253:GLN:HA	1:A:256:GLU:HG2	1.49	0.94
1:A:387:PRO:HA	1:A:390:LYS:HD3	1.50	0.93
1:A:124:ARG:HA	1:A:131:VAL:HG21	1.53	0.91
1:A:363:LYS:HB3	4:A:2112:HOH:O	1.68	0.90
1:A:212:GLN:HA	1:A:215:LYS:HD2	1.60	0.84
1:A:212:GLN:HA	1:A:215:LYS:CD	2.07	0.84
1:A:372:THR:HG21	4:A:2115:HOH:O	1.79	0.83
1:A:153:GLU:HA	1:A:210:PHE:CD2	2.17	0.79
1:A:167:GLN:HE22	1:A:180:ASP:H	1.30	0.79
1:A:363:LYS:CB	4:A:2112:HOH:O	2.28	0.79
1:A:169:SER:O	1:A:173:LEU:HD13	1.82	0.77
1:A:212:GLN:N	1:A:215:LYS:HE3	1.99	0.77
1:A:263:ILE:HD13	1:A:307:TYR:HB2	1.67	0.76
1:A:246:GLY:O	1:A:335:ILE:HD11	1.86	0.75
1:A:361:LEU:HD22	1:A:362:PRO:HD2	1.67	0.75
1:A:44:MSE:O	1:A:48:GLN:HG3	1.85	0.75
1:A:355:SER:OG	1:A:357:LEU:HD13	1.87	0.75
1:A:163:GLN:O	1:A:167:GLN:HG3	1.88	0.74
1:A:253:GLN:HA	1:A:256:GLU:CG	2.18	0.74
1:A:314:GLY:HA3	1:A:318:ASN:ND2	2.02	0.73
1:A:308:PHE:N	1:A:328:ILE:HD11	2.04	0.73
1:A:104:LEU:HB3	1:A:106:ILE:HD11	1.71	0.73
1:A:50:ASN:HB2	1:A:74:TYR:OH	1.91	0.69
1:A:124:ARG:HA	1:A:131:VAL:CG2	2.22	0.68
1:A:30:LEU:HD13	1:A:106:ILE:CD1	2.23	0.67
1:A:385:ASN:HB3	4:A:2118:HOH:O	1.95	0.67
1:A:136:LEU:HB2	1:A:195:VAL:HG12	1.75	0.67
1:A:138:LEU:N	1:A:138:LEU:HD12	2.09	0.66
1:A:314:GLY:O	1:A:317:ILE:HG12	1.96	0.66
1:A:137:ASN:HD22	1:A:198:HIS:HE1	1.44	0.65
1:A:273:ASN:HB2	1:A:276:VAL:HG12	1.79	0.64
1:A:87:PRO:HD2	1:A:119:PRO:HG3	1.80	0.64
1:A:356:SER:HB2	3:A:1415:CSF:H5'	1.80	0.63
1:A:157:ILE:O	1:A:161:ILE:HG12	1.99	0.63
1:A:211:LEU:O	1:A:215:LYS:HG3	2.00	0.62
1:A:328:ILE:HD13	1:A:328:ILE:C	2.20	0.62
1:A:290:HIS:HD2	1:A:297:ASP:H	1.48	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:326:LYS:NZ	1:A:326:LYS:HB2	2.15	0.61
1:A:306:ILE:HD12	1:A:306:ILE:N	2.16	0.60
1:A:328:ILE:HD13	1:A:329:THR:N	2.17	0.60
1:A:300:ILE:HD13	1:A:300:ILE:H	1.68	0.59
1:A:364:GLU:CD	1:A:364:GLU:H	2.05	0.59
1:A:111:ALA:HB1	1:A:144:MSE:HE3	1.85	0.59
1:A:67:PRO:HG2	1:A:70:ILE:CG1	2.32	0.59
1:A:273:ASN:HB2	1:A:276:VAL:CG1	2.33	0.59
1:A:361:LEU:HD12	1:A:366:ILE:CD1	2.32	0.59
1:A:290:HIS:CD2	1:A:297:ASP:H	2.20	0.59
1:A:30:LEU:N	1:A:30:LEU:HD12	2.19	0.58
1:A:402:GLN:C	1:A:402:GLN:HE21	2.07	0.58
1:A:91:LEU:HD21	1:A:120:ILE:HD13	1.84	0.57
1:A:372:THR:CG2	4:A:2115:HOH:O	2.42	0.57
1:A:368:HIS:HB2	4:A:2113:HOH:O	2.05	0.57
1:A:67:PRO:HG3	1:A:243:THR:HG22	1.87	0.57
1:A:361:LEU:HD12	1:A:366:ILE:HD13	1.87	0.56
1:A:346:LEU:HB3	1:A:365:LYS:HE3	1.87	0.56
1:A:309:LYS:CE	1:A:335:ILE:O	2.53	0.56
1:A:163:GLN:HA	1:A:163:GLN:OE1	2.05	0.56
1:A:212:GLN:HA	1:A:215:LYS:CG	2.35	0.55
1:A:274:THR:O	1:A:278:GLU:HG3	2.07	0.55
1:A:285:LEU:C	1:A:285:LEU:HD13	2.27	0.55
1:A:190:GLN:NE2	1:A:221:ASN:HD22	2.05	0.54
1:A:87:PRO:HD2	1:A:119:PRO:CG	2.37	0.54
1:A:224:LYS:CB	4:A:2066:HOH:O	2.54	0.54
1:A:289:ASN:HB3	1:A:293:GLN:HG3	1.90	0.54
1:A:306:ILE:HD12	1:A:306:ILE:H	1.72	0.54
1:A:67:PRO:HG2	1:A:70:ILE:HG12	1.90	0.54
1:A:89:GLU:HG2	1:A:92:PHE:CE2	2.43	0.54
1:A:33:ASP:CG	1:A:38:PRO:HB2	2.28	0.53
1:A:399:ASP:OD1	1:A:401:SER:HB3	2.07	0.53
1:A:67:PRO:HA	4:A:2019:HOH:O	2.06	0.53
1:A:128:LEU:HD22	1:A:128:LEU:H	1.74	0.52
1:A:326:LYS:HB2	1:A:326:LYS:HZ2	1.74	0.52
1:A:137:ASN:C	1:A:138:LEU:HD12	2.29	0.52
1:A:293:GLN:HB3	1:A:295:GLU:OE1	2.10	0.52
1:A:224:LYS:HB3	4:A:2066:HOH:O	2.08	0.52
1:A:393:ARG:NH2	1:A:403:VAL:HG23	2.26	0.51
1:A:167:GLN:NE2	1:A:180:ASP:H	2.03	0.51
1:A:224:LYS:HD2	1:A:225:MSE:H	1.74	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:100:GLY:O	1:A:130:ARG:HD3	2.11	0.51
1:A:393:ARG:NH2	1:A:401:SER:O	2.44	0.50
1:A:152:GLU:OE2	1:A:155:LYS:HD2	2.11	0.50
1:A:85:ASN:ND2	1:A:112:HIS:CG	2.80	0.50
1:A:392:MSE:HE2	1:A:397:ILE:HD12	1.93	0.50
1:A:328:ILE:HD13	1:A:329:THR:C	2.32	0.49
1:A:92:PHE:CE1	1:A:123:TYR:HB2	2.47	0.49
1:A:136:LEU:HB2	1:A:195:VAL:CG1	2.42	0.48
1:A:260:ALA:HA	1:A:348:ASP:OD2	2.13	0.48
1:A:212:GLN:HB2	1:A:213:PRO:HD3	1.95	0.48
1:A:37:LEU:HB2	1:A:38:PRO:HD3	1.96	0.48
1:A:212:GLN:HA	1:A:215:LYS:HG3	1.95	0.48
1:A:91:LEU:O	1:A:95:LEU:HD22	2.13	0.48
1:A:37:LEU:HD22	1:A:37:LEU:H	1.79	0.47
1:A:90:ALA:O	1:A:94:ILE:HG12	2.14	0.47
1:A:118:ARG:HB3	1:A:119:PRO:HD3	1.96	0.47
1:A:136:LEU:HD12	1:A:195:VAL:HG11	1.96	0.47
1:A:304:TYR:HB2	4:A:2093:HOH:O	2.13	0.47
1:A:316:GLU:HG3	1:A:317:ILE:N	2.29	0.47
1:A:224:LYS:HD2	1:A:225:MSE:N	2.30	0.47
1:A:83:LYS:O	1:A:86:ARG:HG2	2.15	0.47
1:A:138:LEU:HD22	1:A:197:TYR:CE1	2.49	0.47
1:A:153:GLU:HA	1:A:210:PHE:CE2	2.48	0.47
1:A:240:PHE:O	1:A:243:THR:HB	2.15	0.47
1:A:290:HIS:HE1	1:A:410:LYS:O	1.98	0.47
1:A:103:GLU:HG3	4:A:2003:HOH:O	2.15	0.46
1:A:308:PHE:CB	1:A:328:ILE:HD11	2.45	0.46
1:A:253:GLN:CA	1:A:256:GLU:HG2	2.34	0.46
1:A:307:TYR:CD2	1:A:331:ILE:HD11	2.51	0.46
1:A:314:GLY:HA3	1:A:318:ASN:HD21	1.80	0.45
1:A:138:LEU:N	1:A:138:LEU:CD1	2.77	0.45
1:A:340:LEU:N	1:A:340:LEU:HD23	2.31	0.45
1:A:26:LYS:HB2	1:A:102:ILE:HG22	1.99	0.45
1:A:118:ARG:HB3	1:A:119:PRO:CD	2.47	0.45
1:A:261:LYS:HE2	4:A:2083:HOH:O	2.17	0.44
1:A:51:GLU:O	1:A:53:LYS:HD3	2.17	0.44
1:A:70:ILE:HA	1:A:73:GLN:HE21	1.82	0.44
1:A:300:ILE:HG12	1:A:304:TYR:CD2	2.52	0.44
1:A:355:SER:OG	1:A:357:LEU:CD1	2.62	0.44
1:A:406:TRP:HA	1:A:406:TRP:HE3	1.82	0.44
1:A:78:HIS:HD2	4:A:2012:HOH:O	1.99	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:258:GLN:CD	1:A:258:GLN:C	2.77	0.44
1:A:285:LEU:O	1:A:288:LEU:HB3	2.18	0.44
1:A:291:PHE:O	1:A:299:PHE:HA	2.18	0.44
1:A:385:ASN:CB	4:A:2118:HOH:O	2.61	0.44
1:A:101:ASN:HD22	1:A:101:ASN:HA	1.59	0.44
1:A:227:TRP:CH2	1:A:338:GLU:HG2	2.52	0.44
1:A:273:ASN:CB	4:A:2089:HOH:O	2.16	0.44
1:A:313:ARG:HD3	4:A:2096:HOH:O	2.17	0.44
1:A:200:LEU:HD22	1:A:359:PHE:HB3	2.00	0.44
1:A:101:ASN:HA	1:A:130:ARG:HG2	1.99	0.44
1:A:314:GLY:O	1:A:317:ILE:CG1	2.63	0.44
1:A:111:ALA:HB1	1:A:144:MSE:CE	2.47	0.43
1:A:273:ASN:N	1:A:273:ASN:HD22	2.15	0.43
1:A:214:LEU:HD12	1:A:218:LEU:HG	2.00	0.43
1:A:313:ARG:NH1	4:A:2096:HOH:O	2.48	0.43
1:A:237:GLN:O	1:A:240:PHE:HB3	2.19	0.43
1:A:406:TRP:HA	1:A:406:TRP:CE3	2.53	0.43
1:A:355:SER:HA	4:A:2129:HOH:O	2.18	0.43
1:A:36:SER:HB2	4:A:2005:HOH:O	2.19	0.43
1:A:248:ASN:OD1	1:A:250:GLU:HB3	2.19	0.42
1:A:326:LYS:HG2	1:A:327:ASN:N	2.34	0.42
1:A:307:TYR:HD2	1:A:331:ILE:HD11	1.83	0.42
1:A:66:ILE:HD13	1:A:79:PHE:HE2	1.84	0.42
1:A:208:ALA:HB1	1:A:210:PHE:CE1	2.54	0.42
1:A:307:TYR:HA	1:A:329:THR:O	2.19	0.42
1:A:31:TYR:CD1	1:A:42:GLN:HB3	2.54	0.42
1:A:308:PHE:HB2	1:A:328:ILE:HD11	2.02	0.42
1:A:309:LYS:HE3	1:A:335:ILE:O	2.19	0.42
1:A:85:ASN:HD22	1:A:112:HIS:CG	2.37	0.42
1:A:51:GLU:O	1:A:53:LYS:CD	2.68	0.41
1:A:289:ASN:HB3	1:A:293:GLN:CG	2.49	0.41
1:A:361:LEU:HD22	1:A:362:PRO:CD	2.43	0.41
1:A:50:ASN:HD22	1:A:51:GLU:N	2.19	0.41
1:A:357:LEU:HD22	1:A:358:TYR:CD1	2.56	0.41
1:A:210:PHE:CE2	1:A:211:LEU:HD13	2.56	0.41
1:A:85:ASN:ND2	1:A:112:HIS:CD2	2.88	0.41
1:A:125:PHE:O	1:A:128:LEU:HD22	2.21	0.41
1:A:127:HIS:O	1:A:131:VAL:HG22	2.20	0.41
1:A:30:LEU:HD13	1:A:106:ILE:HD13	2.00	0.41
1:A:86:ARG:HA	1:A:87:PRO:HD3	1.93	0.40
1:A:247:PHE:HZ	1:A:345:LEU:HD12	1.86	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:294:ALA:HA	1:A:299:PHE:CD2	2.57	0.40
1:A:361:LEU:HD12	1:A:366:ILE:HD11	2.02	0.40
1:A:171:TYR:HA	1:A:176:LYS:O	2.22	0.40
1:A:110:ILE:HG12	1:A:138:LEU:HB3	2.03	0.40
1:A:356:SER:HB2	3:A:1415:CSF:C5'	2.50	0.40

All (2) 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)
4:A:2027:HOH:O	4:A:2045:HOH:O 4_456	1.80	0.40
4:A:2078:HOH:O	4:A:2112:HOH:O 3_655	1.83	0.37

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	373/389 (96%)	341 (91%)	28 (8%)	4 (1%)	14 26

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	401	SER
1	A	406	TRP
1	A	219	ALA
1	A	332	PRO

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	341/346 (99%)	322 (94%)	19 (6%)	21	40

All (19) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	50	ASN
1	A	95	LEU
1	A	124	ARG
1	A	163	GLN
1	A	200	LEU
1	A	211	LEU
1	A	242	LEU
1	A	258	GLN
1	A	276	VAL
1	A	295	GLU
1	A	300	ILE
1	A	302	ASP
1	A	328	ILE
1	A	346	LEU
1	A	357	LEU
1	A	368	HIS
1	A	388	TYR
1	A	402	GLN
1	A	406	TRP

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (17) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	50	ASN
1	A	55	HIS
1	A	73	GLN
1	A	85	ASN
1	A	101	ASN
1	A	112	HIS
1	A	137	ASN
1	A	154	ASN
1	A	167	GLN
1	A	190	GLN

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Mol	Chain	Res	Type
1	A	212	GLN
1	A	238	GLN
1	A	253	GLN
1	A	273	ASN
1	A	282	GLN
1	A	290	HIS
1	A	402	GLN

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](#)

2 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	BGC	B	1	2	12,12,12	1.61	4 (33%)	17,17,17	0.84	0
2	GAL	B	2	2	11,11,12	2.00	4 (36%)	15,15,17	1.03	1 (6%)

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	BGC	B	1	2	-	0/2/22/22	0/1/1/1
2	GAL	B	2	2	-	0/2/19/22	0/1/1/1

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	2	GAL	C6-C5	-4.64	1.36	1.51
2	B	1	BGC	O5-C1	2.84	1.50	1.42
2	B	1	BGC	C1-C2	2.62	1.58	1.52
2	B	2	GAL	C2-C3	2.51	1.56	1.52
2	B	1	BGC	O5-C5	2.25	1.49	1.44
2	B	2	GAL	C4-C3	2.05	1.57	1.52
2	B	2	GAL	O5-C5	2.03	1.47	1.43
2	B	1	BGC	C4-C3	2.01	1.57	1.52

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	2	GAL	O5-C5-C6	2.16	110.58	107.20

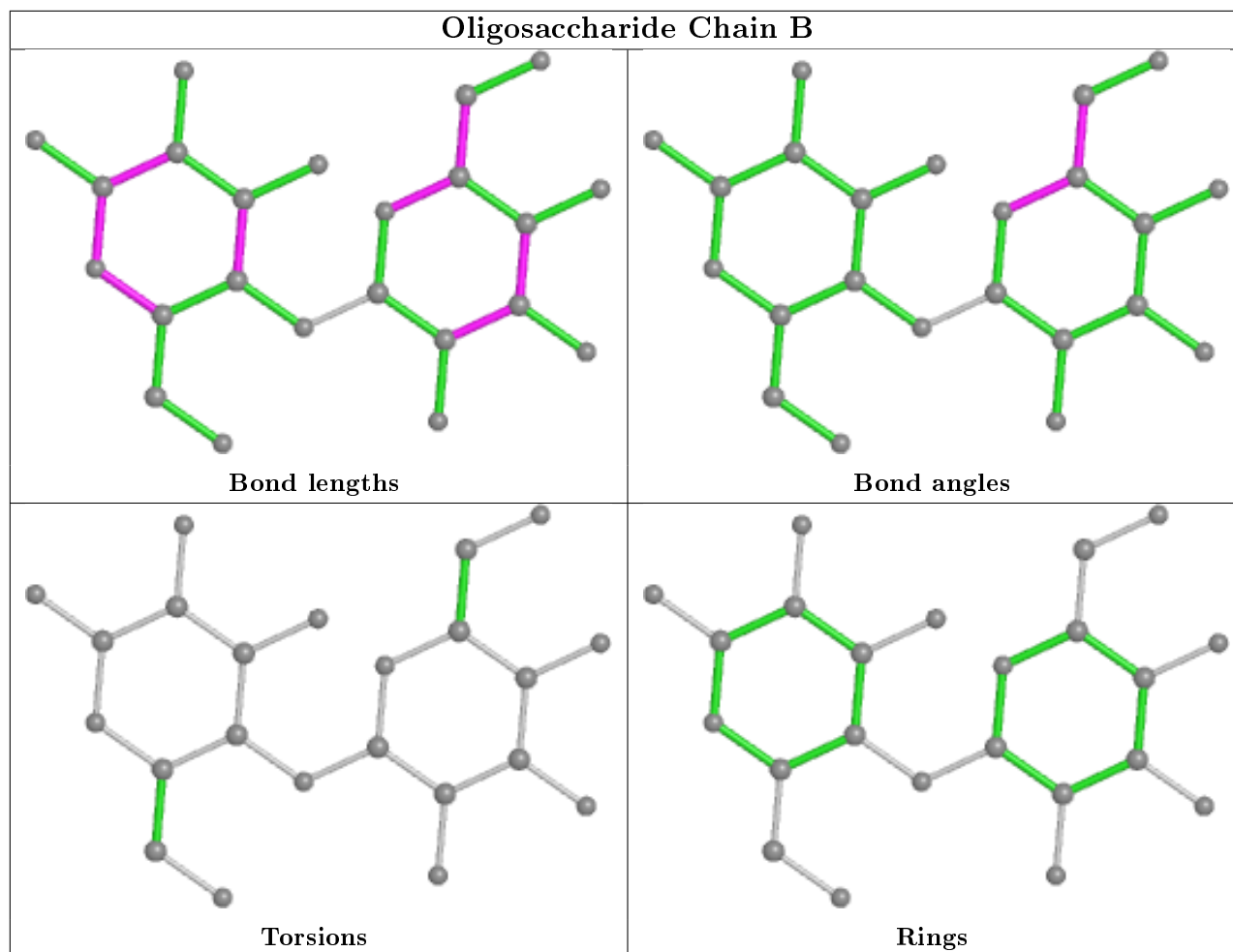
There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

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



5.6 Ligand geometry [i](#)

1 ligand is 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	CSF	A	1415	-	36,44,44	1.98	11 (30%)	42,67,67	3.03	20 (47%)

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	CSF	A	1415	-	1/1/14/15	10/25/75/75	0/3/3/3

All (11) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	1415	CSF	C4-N4	-4.96	1.20	1.35
3	A	1415	CSF	PA-O2A	-3.78	1.37	1.50
3	A	1415	CSF	O10-C10	-3.60	1.15	1.23
3	A	1415	CSF	PA-O3A	-3.34	1.39	1.55
3	A	1415	CSF	C5A-N5A	-2.79	1.41	1.45
3	A	1415	CSF	O4'-C4'	-2.66	1.39	1.45
3	A	1415	CSF	C6-N1	2.49	1.38	1.35
3	A	1415	CSF	F3A-C3A	-2.24	1.36	1.40
3	A	1415	CSF	C2'-C1'	-2.24	1.50	1.53
3	A	1415	CSF	C3A-C4A	2.16	1.54	1.52
3	A	1415	CSF	O6A-C6A	-2.01	1.40	1.44

All (20) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	1415	CSF	N4-C4-N3	10.12	132.49	116.49
3	A	1415	CSF	C5-C4-N4	-8.12	107.04	121.14
3	A	1415	CSF	O7A-C7A-C8A	-5.71	95.03	108.81
3	A	1415	CSF	C2-N3-C4	5.13	121.54	116.34
3	A	1415	CSF	O4A-C4A-C5A	3.90	117.53	109.66
3	A	1415	CSF	O7A-C7A-C6A	-3.68	101.55	109.50
3	A	1415	CSF	O10-C10-C11	-3.53	115.51	122.06
3	A	1415	CSF	O1A-PA-O2A	3.43	121.87	109.29
3	A	1415	CSF	O8A-C8A-C9A	-3.40	101.18	109.14
3	A	1415	CSF	C4A-C5A-N5A	3.34	116.93	110.62
3	A	1415	CSF	O3'-C3'-C4'	2.88	119.36	111.05
3	A	1415	CSF	C3A-C4A-C5A	-2.74	106.30	109.83
3	A	1415	CSF	C5A-N5A-C10	2.72	129.80	123.18
3	A	1415	CSF	C8A-C7A-C6A	2.67	118.10	113.03
3	A	1415	CSF	O8A-C8A-C7A	-2.66	102.63	109.10
3	A	1415	CSF	C6-N1-C2	-2.66	116.97	121.20
3	A	1415	CSF	O4'-C1'-C2'	-2.47	103.32	106.93
3	A	1415	CSF	F3A-C3A-C4A	2.37	110.96	108.85
3	A	1415	CSF	C2'-C3'-C4'	-2.09	98.58	102.64
3	A	1415	CSF	O10-C10-N5A	2.05	125.72	121.95

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
3	A	1415	CSF	C7A

All (10) torsion outliers are listed below:

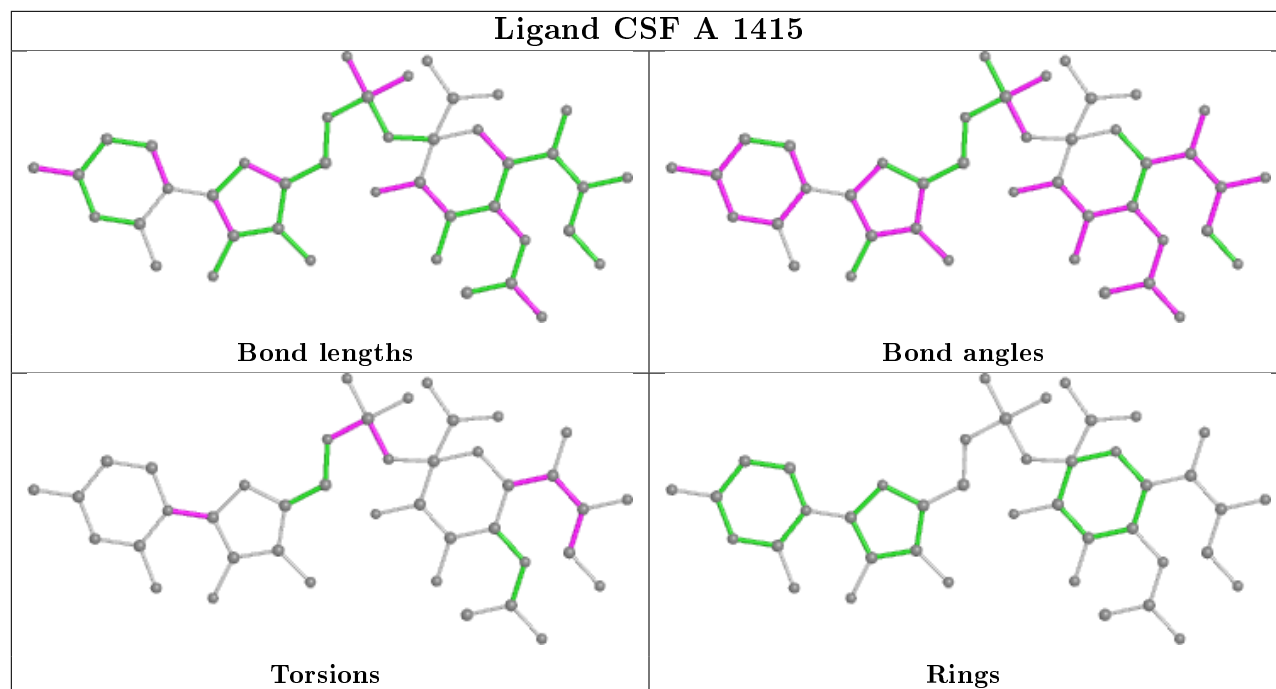
Mol	Chain	Res	Type	Atoms
3	A	1415	CSF	C5A-C6A-C7A-O7A
3	A	1415	CSF	C6A-C7A-C8A-C9A
3	A	1415	CSF	C2A-O1A-PA-O5'
3	A	1415	CSF	O4'-C1'-N1-C6
3	A	1415	CSF	C5'-O5'-PA-O1A
3	A	1415	CSF	C5'-O5'-PA-O3A
3	A	1415	CSF	C6A-C7A-C8A-O8A
3	A	1415	CSF	O7A-C7A-C8A-C9A
3	A	1415	CSF	C7A-C8A-C9A-O9A
3	A	1415	CSF	C5'-O5'-PA-O2A

There are no ring outliers.

1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	1415	CSF	2	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.



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	371/389 (95%)	-0.52	1 (0%) 94 94	4, 16, 30, 39	0

All (1) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	125	PHE	3.1

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

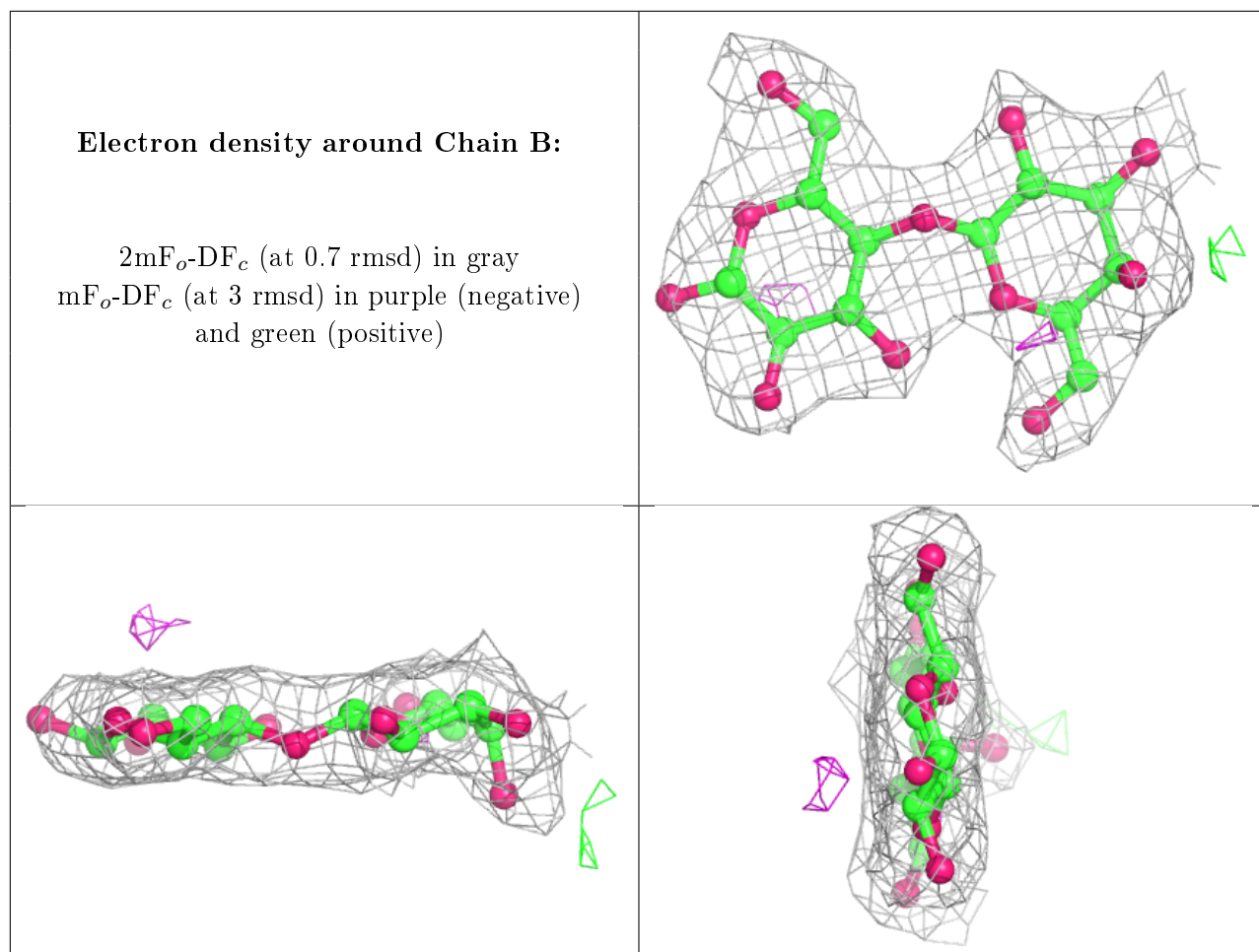
There are no non-standard protein/DNA/RNA residues in this entry.

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	GAL	B	2	11/12	0.89	0.17	21,22,24,27	0
2	BGC	B	1	12/12	0.94	0.15	23,24,27,29	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.

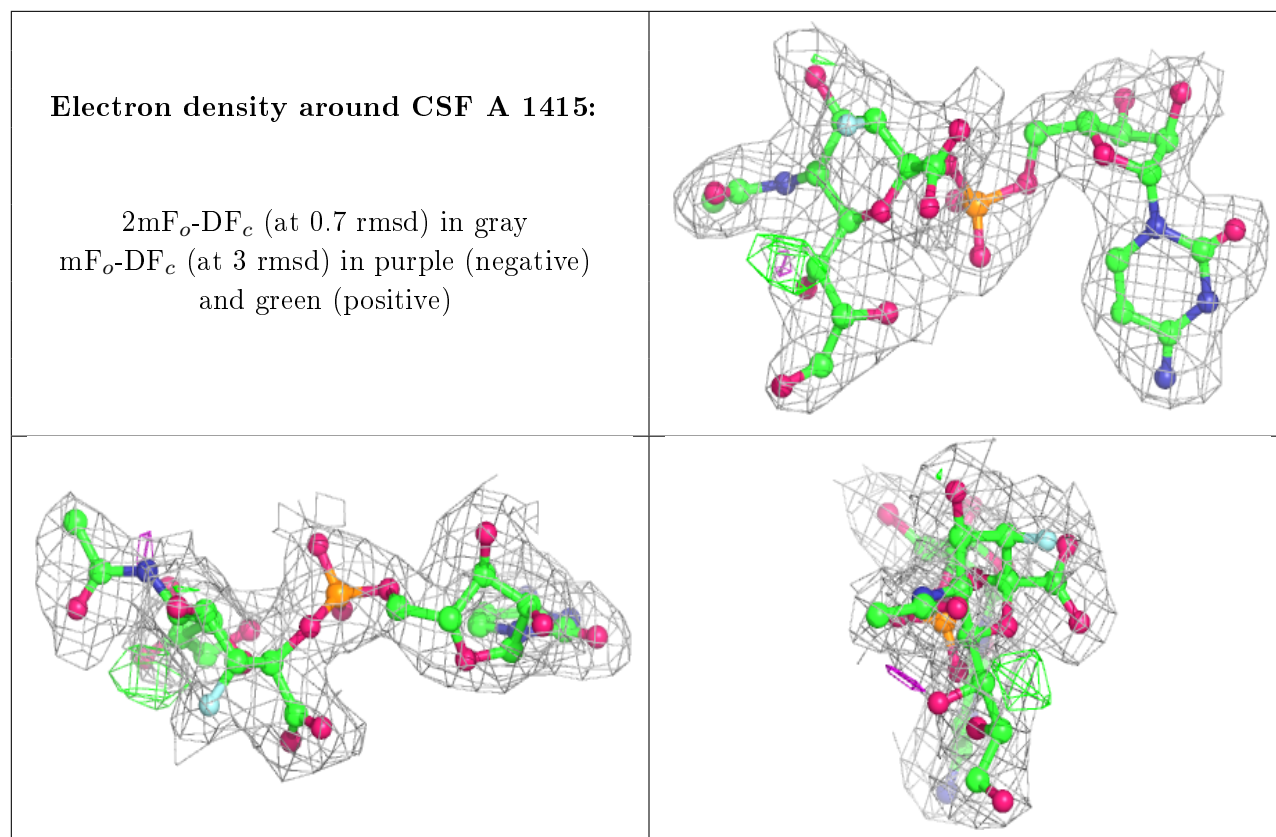


6.4 Ligands [i](#)

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
3	CSF	A	1415	42/42	0.95	0.12	4,26,37,41	0

The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.



6.5 Other polymers [i](#)

There are no such residues in this entry.