



Full wwPDB X-ray Structure Validation Report ⓘ

Aug 9, 2020 – 04:12 PM BST

PDB ID : 3ESW
Title : Complex of yeast PNGase with GlcNAc2-IAc.
Authors : Zhao, G.; Zhou, X.; Lennarz, W.J.; Schindelin, H.
Deposited on : 2008-10-06
Resolution : 3.40 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
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

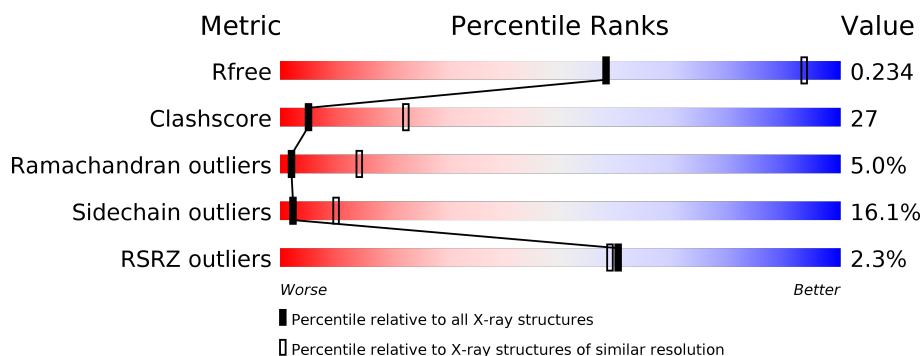
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.40 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	1026 (3.48-3.32)
Clashscore	141614	1055 (3.48-3.32)
Ramachandran outliers	138981	1038 (3.48-3.32)
Sidechain outliers	138945	1038 (3.48-3.32)
RSRZ outliers	127900	2173 (3.50-3.30)

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 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	355	<div> <div>2%</div> <div> <div></div> <div>52%</div> <div>29%</div> <div>11%</div> <div>6%</div> </div> </div>
2	B	55	<div> <div>4%</div> <div> <div></div> <div>29%</div> <div>55%</div> <div>13%</div> <div></div> </div> </div>
3	C	2	<div> <div></div> <div>100%</div> </div>

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

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
1	YCM	A	191	-	-	X	-
3	NAG	C	1	-	-	X	-
4	ZN	A	344	-	-	X	-

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 3220 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 Peptide-N(4)-(N-acetyl-beta-glucosaminyl)asparagine amidase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	333	Total	C	N	O	S	0	0	0
			2761	1755	476	512	18			

There are 21 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-13	MET	-	expression tag	UNP Q02890
A	-12	GLY	-	expression tag	UNP Q02890
A	-11	SER	-	expression tag	UNP Q02890
A	-10	SER	-	expression tag	UNP Q02890
A	-9	HIS	-	expression tag	UNP Q02890
A	-8	HIS	-	expression tag	UNP Q02890
A	-7	HIS	-	expression tag	UNP Q02890
A	-6	HIS	-	expression tag	UNP Q02890
A	-5	HIS	-	expression tag	UNP Q02890
A	-4	HIS	-	expression tag	UNP Q02890
A	-3	SER	-	expression tag	UNP Q02890
A	-2	SER	-	expression tag	UNP Q02890
A	-1	GLY	-	expression tag	UNP Q02890
A	0	LEU	-	expression tag	UNP Q02890
A	1	VAL	-	expression tag	UNP Q02890
A	2	PRO	-	expression tag	UNP Q02890
A	3	ARG	-	expression tag	UNP Q02890
A	4	GLY	-	expression tag	UNP Q02890
A	5	SER	-	expression tag	UNP Q02890
A	6	HIS	-	expression tag	UNP Q02890
A	7	MET	-	expression tag	UNP Q02890

- Molecule 2 is a protein called UV excision repair protein RAD23.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	55	Total	C	N	O	S	0	0	0
			423	269	71	81	2			

- Molecule 3 is an oligosaccharide called 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
3	C	2	Total	C	N	O	0	0	0
			28	16	2	10			

- Molecule 4 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	1	Total	Zn	0	0
			1	1		

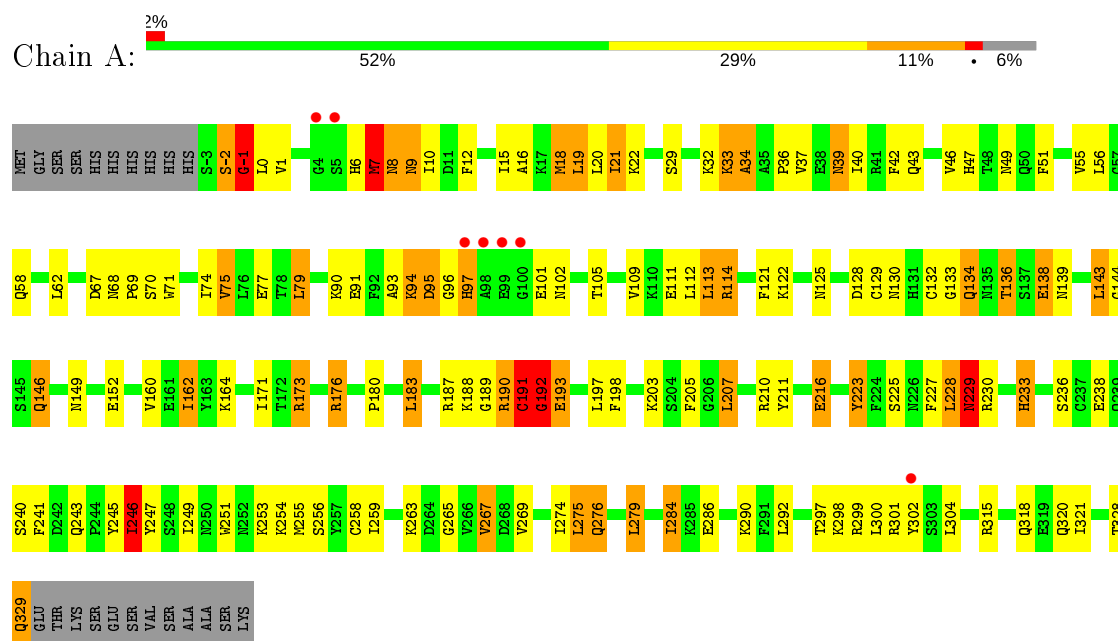
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	7	Total	O	0	0
			7	7		

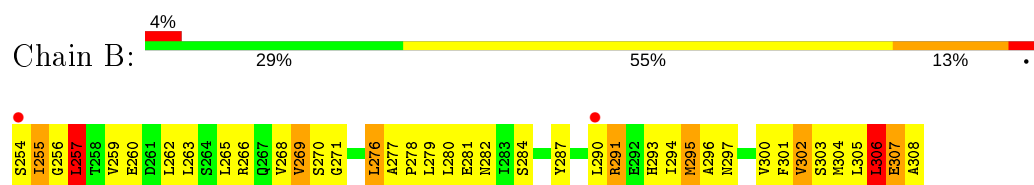
3 Residue-property plots [i](#)

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.

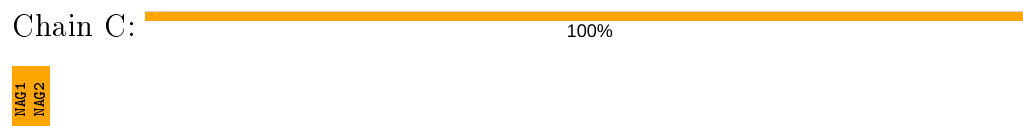
- Molecule 1: Peptide-N(4)-(N-acetyl-beta-glucosaminyl)asparagine amidase



- Molecule 2: UV excision repair protein RAD23



- Molecule 3: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose



4 Data and refinement statistics

Property	Value	Source
Space group	P 31 2 1	Depositor
Cell constants a, b, c, α , β , γ	131.53Å 131.53Å 127.75Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	37.96 – 3.40 37.97 – 3.40	Depositor EDS
% Data completeness (in resolution range)	100.0 (37.96-3.40) 100.0 (37.97-3.40)	Depositor EDS
R_{merge}	0.13	Depositor
R_{sym}	0.13	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.01 (at 3.40Å)	Xtriage
Refinement program	REFMAC 5.5.0054	Depositor
R, R_{free}	0.197 , 0.235 0.196 , 0.234	Depositor DCC
R_{free} test set	914 reflections (5.09%)	wwPDB-VP
Wilson B-factor (Å ²)	75.1	Xtriage
Anisotropy	0.810	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 88.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	0.045 for -h,-k,l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	3220	wwPDB-VP
Average B, all atoms (Å ²)	63.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.35% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: ZN, NAG, YCM

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	0.80	0/2816	0.79	2/3799 (0.1%)
2	B	0.68	0/429	0.85	2/584 (0.3%)
All	All	0.79	0/3245	0.80	4/4383 (0.1%)

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

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	5

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
2	B	306	LEU	CA-CB-CG	6.22	129.61	115.30
1	A	-1	GLY	N-CA-C	5.70	127.36	113.10
2	B	257	LEU	CA-CB-CG	5.47	127.89	115.30
1	A	192	GLY	N-CA-C	5.07	125.76	113.10

There are no chirality outliers.

All (5) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	-1	GLY	Peptide
1	A	-2	SER	Peptide
1	A	190	ARG	Mainchain
1	A	191	YCM	Peptide

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Mol	Chain	Res	Type	Group
1	A	228	LEU	Peptide

5.2 Too-close contacts ⓘ

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2761	0	2669	123	0
2	B	423	0	434	57	0
3	C	28	0	26	10	0
4	A	1	0	0	2	0
5	A	7	0	0	0	0
All	All	3220	0	3129	170	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 27.

All (170) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:265:LEU:CD2	2:B:305:LEU:HD21	1.71	1.20
2:B:265:LEU:HD21	2:B:305:LEU:HD21	1.16	1.07
1:A:20:LEU:HD11	2:B:276:LEU:HD23	1.37	1.06
1:A:19:LEU:HD23	2:B:276:LEU:HD21	1.46	0.98
2:B:255:ILE:HD13	2:B:256:GLY:N	1.84	0.91
1:A:134:GLN:O	1:A:136:THR:HG22	1.70	0.90
1:A:191:YCM:CE	3:C:1:NAG:C1	2.49	0.90
2:B:255:ILE:CD1	2:B:257:LEU:HD22	2.01	0.90
3:C:1:NAG:C4	3:C:2:NAG:C1	2.48	0.90
1:A:20:LEU:CD1	2:B:276:LEU:HD23	2.03	0.88
2:B:265:LEU:HD21	2:B:305:LEU:CD2	2.03	0.86
2:B:276:LEU:HD12	2:B:276:LEU:O	1.77	0.85
2:B:255:ILE:HD12	2:B:257:LEU:HD22	1.58	0.84
1:A:176:ARG:HG2	1:A:176:ARG:HH11	1.43	0.84
1:A:15:ILE:O	1:A:18:MET:O	1.96	0.83
2:B:303:SER:O	2:B:306:LEU:HD13	1.81	0.81
1:A:143:LEU:HD12	1:A:143:LEU:N	1.97	0.79

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:265:LEU:HD12	2:B:279:LEU:HD23	1.70	0.73
1:A:36:PRO:O	1:A:40:ILE:HD12	1.88	0.72
1:A:12:PHE:HA	1:A:15:ILE:HG22	1.72	0.70
1:A:152:GLU:OE1	1:A:173:ARG:NH1	2.26	0.69
1:A:36:PRO:HA	1:A:39:ASN:HB2	1.75	0.68
1:A:144:GLY:O	1:A:162:ILE:HD13	1.93	0.68
1:A:251:TRP:CZ3	3:C:1:NAG:H83	2.29	0.67
1:A:233:HIS:O	1:A:241:PHE:O	2.13	0.67
1:A:180:PRO:HA	1:A:183:LEU:HD22	1.76	0.67
1:A:113:LEU:HD13	1:A:198:PHE:CZ	2.30	0.66
2:B:265:LEU:CD2	2:B:305:LEU:CD2	2.62	0.66
2:B:255:ILE:HD12	2:B:257:LEU:CD2	2.26	0.65
2:B:269:VAL:HG22	2:B:269:VAL:O	1.96	0.65
1:A:176:ARG:CG	1:A:176:ARG:HH11	2.10	0.65
1:A:230:ARG:HD3	1:A:328:THR:HG23	1.78	0.64
1:A:143:LEU:CD1	1:A:143:LEU:N	2.59	0.64
1:A:191:YCM:O	1:A:192:GLY:C	2.35	0.64
1:A:7:MET:HE3	1:A:8:ASN:O	1.98	0.63
2:B:306:LEU:CD2	2:B:306:LEU:O	2.46	0.63
1:A:10:ILE:HD11	2:B:294:ILE:HG22	1.82	0.62
1:A:15:ILE:HD12	1:A:18:MET:HG3	1.82	0.62
2:B:303:SER:HA	2:B:306:LEU:HD12	1.82	0.61
1:A:205:PHE:HB2	1:A:207:LEU:HD22	1.82	0.61
1:A:96:GLY:O	1:A:97:HIS:O	2.18	0.61
3:C:1:NAG:H4	3:C:2:NAG:C1	2.32	0.60
1:A:132:CYS:HG	4:A:344:ZN:ZN	1.16	0.60
1:A:132:CYS:SG	4:A:344:ZN:ZN	1.89	0.59
1:A:246:ILE:HG22	1:A:247:TYR:N	2.17	0.59
1:A:251:TRP:HZ3	3:C:1:NAG:H83	1.66	0.59
1:A:152:GLU:OE1	1:A:160:VAL:HG21	2.02	0.59
1:A:125:ASN:O	1:A:190:ARG:NH1	2.36	0.59
1:A:18:MET:SD	1:A:19:LEU:N	2.75	0.59
1:A:210:ARG:NH2	1:A:267:VAL:HG11	2.17	0.59
1:A:19:LEU:HD13	2:B:302:VAL:HG22	1.85	0.59
2:B:306:LEU:HD22	2:B:307:GLU:OE1	2.03	0.58
2:B:306:LEU:O	2:B:306:LEU:HD22	2.04	0.58
1:A:328:THR:HG22	1:A:329:GLN:H	1.69	0.58
1:A:263:LYS:O	1:A:300:LEU:HD22	2.05	0.57
2:B:303:SER:HA	2:B:306:LEU:CD1	2.34	0.57
1:A:6:HIS:CD2	1:A:9:ASN:HB2	2.41	0.56
1:A:7:MET:HE3	1:A:8:ASN:HB3	1.86	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:129:CYS:H	1:A:136:THR:CG2	2.19	0.55
1:A:129:CYS:H	1:A:136:THR:HG23	1.71	0.55
1:A:247:TYR:HB3	1:A:255:MET:HE3	1.88	0.55
1:A:46:VAL:O	1:A:46:VAL:HG12	2.05	0.55
1:A:138:GLU:H	1:A:138:GLU:CD	2.10	0.55
2:B:262:LEU:HD13	2:B:308:ALA:CB	2.37	0.55
2:B:276:LEU:C	2:B:276:LEU:HD12	2.26	0.55
2:B:297:ASN:HB3	2:B:300:VAL:HG12	1.87	0.55
1:A:71:TRP:HA	1:A:74:ILE:HD12	1.89	0.54
1:A:139:ASN:OD1	1:A:139:ASN:N	2.41	0.54
1:A:284:ILE:HG23	1:A:284:ILE:O	2.07	0.54
1:A:16:ALA:O	1:A:20:LEU:HD13	2.08	0.54
1:A:7:MET:HE3	1:A:8:ASN:CB	2.38	0.54
1:A:34:ALA:HB2	1:A:302:TYR:OH	2.08	0.54
1:A:51:PHE:O	1:A:55:VAL:HG23	2.08	0.54
1:A:105:THR:HG22	1:A:223:TYR:OH	2.09	0.53
1:A:2:SER:HA	1:A:0:LEU:H	1.74	0.53
1:A:191:YCM:NZ2	3:C:1:NAG:O5	2.39	0.53
1:A:227:PHE:CD2	1:A:228:LEU:HD23	2.44	0.53
1:A:29:SER:O	1:A:32:LYS:HB2	2.08	0.53
1:A:7:MET:CE	2:B:291:ARG:HH12	2.22	0.52
1:A:18:MET:O	1:A:19:LEU:HB3	2.10	0.52
1:A:267:VAL:HG22	1:A:269:VAL:HG13	1.92	0.52
1:A:10:ILE:CD1	2:B:294:ILE:HG22	2.39	0.51
1:A:90:LYS:O	1:A:93:ALA:HB3	2.11	0.51
1:A:19:LEU:HD13	2:B:302:VAL:CG2	2.40	0.51
2:B:262:LEU:HD13	2:B:308:ALA:HB1	1.93	0.51
1:A:136:THR:O	1:A:136:THR:HG23	2.11	0.51
1:A:162:ILE:H	1:A:162:ILE:HD13	1.75	0.51
1:A:171:ILE:HD12	1:A:171:ILE:N	2.26	0.51
2:B:255:ILE:HD11	2:B:257:LEU:HD22	1.91	0.50
2:B:268:VAL:C	2:B:270:SER:H	2.14	0.50
1:A:113:LEU:HD12	1:A:236:SER:HA	1.93	0.50
1:A:228:LEU:O	1:A:230:ARG:N	2.44	0.50
1:A:19:LEU:CD2	2:B:276:LEU:HD21	2.29	0.50
2:B:276:LEU:HD11	2:B:280:LEU:HD11	1.94	0.50
2:B:277:ALA:HB3	2:B:278:PRO:CD	2.42	0.49
1:A:19:LEU:CD1	2:B:302:VAL:HG22	2.41	0.49
1:A:301:ARG:HB3	1:A:304:LEU:HD12	1.95	0.49
1:A:265:GLY:HA3	1:A:297:THR:HG23	1.94	0.48
2:B:255:ILE:HG12	2:B:290:LEU:HD11	1.95	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:58:GLN:HB3	1:A:62:LEU:HD12	1.95	0.48
2:B:284:SER:HB2	2:B:291:ARG:HG2	1.96	0.48
1:A:18:MET:O	1:A:19:LEU:CB	2.61	0.48
2:B:295:MET:CE	2:B:296:ALA:HB2	2.43	0.48
1:A:216:GLU:HG3	1:A:255:MET:HA	1.94	0.48
1:A:75:VAL:HG23	1:A:79:LEU:HD22	1.96	0.47
1:A:225:SER:O	1:A:229:ASN:HA	2.14	0.47
1:A:109:VAL:HA	1:A:112:LEU:HD12	1.95	0.47
1:A:21:ILE:HA	1:A:21:ILE:HD13	1.79	0.47
2:B:259:VAL:HG23	2:B:260:GLU:N	2.29	0.47
1:A:321:ILE:HD11	2:B:266:ARG:HE	1.80	0.47
2:B:287:TYR:HB3	2:B:290:LEU:HD23	1.97	0.47
2:B:255:ILE:HD13	2:B:256:GLY:CA	2.45	0.46
2:B:278:PRO:O	2:B:281:GLU:HB2	2.15	0.46
1:A:7:MET:HE2	2:B:291:ARG:HH12	1.79	0.46
1:A:227:PHE:CD2	1:A:228:LEU:CD2	2.99	0.46
1:A:256:SER:O	1:A:274:ILE:HD12	2.15	0.46
2:B:277:ALA:HB3	2:B:278:PRO:HD3	1.98	0.46
1:A:191:YCM:SG	1:A:192:GLY:N	2.89	0.46
1:A:146:GLN:HG2	1:A:162:ILE:HD12	1.97	0.46
1:A:7:MET:CE	1:A:8:ASN:O	2.63	0.46
1:A:128:ASP:CG	1:A:134:GLN:O	2.54	0.45
1:A:176:ARG:NH1	1:A:176:ARG:HG2	2.20	0.45
2:B:257:LEU:HD23	2:B:257:LEU:H	1.81	0.45
2:B:305:LEU:HD23	2:B:305:LEU:O	2.16	0.45
2:B:254:SER:HB3	2:B:255:ILE:HA	1.98	0.45
1:A:149:ASN:OD1	1:A:149:ASN:C	2.54	0.45
1:A:111:GLU:OE1	1:A:114:ARG:NH1	2.49	0.45
2:B:265:LEU:HD12	2:B:279:LEU:CD2	2.45	0.45
3:C:1:NAG:O4	3:C:2:NAG:C2	2.57	0.45
1:A:253:LYS:HD2	3:C:1:NAG:H81	1.98	0.44
1:A:43:GLN:O	1:A:47:HIS:ND1	2.48	0.44
1:A:67:ASP:OD2	1:A:203:LYS:NZ	2.43	0.44
1:A:243:GLN:O	1:A:246:ILE:HB	2.18	0.44
1:A:101:GLU:O	1:A:102:ASN:HB2	2.18	0.44
1:A:19:LEU:O	1:A:22:LYS:N	2.51	0.44
1:A:40:ILE:H	1:A:40:ILE:HD12	1.83	0.44
2:B:305:LEU:C	2:B:307:GLU:H	2.21	0.44
1:A:189:GLY:HA3	1:A:193:GLU:HG3	2.00	0.44
1:A:299:ARG:HH11	1:A:299:ARG:HG3	1.81	0.44
1:A:33:LYS:O	1:A:34:ALA:CB	2.66	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:191:YCM:NZ2	3:C:1:NAG:C2	2.68	0.43
1:A:238:GLU:OE1	3:C:2:NAG:N2	2.40	0.43
1:A:275:LEU:O	1:A:276:GLN:HB3	2.18	0.43
1:A:328:THR:HG22	1:A:329:GLN:N	2.32	0.43
1:A:246:ILE:CG2	1:A:247:TYR:N	2.80	0.43
1:A:143:LEU:HD11	1:A:164:LYS:CB	2.48	0.43
1:A:122:LYS:O	1:A:188:LYS:HA	2.17	0.43
1:A:255:MET:N	1:A:279:LEU:HD12	2.33	0.43
1:A:143:LEU:HD11	1:A:164:LYS:HB2	2.00	0.43
1:A:62:LEU:HD22	1:A:211:TYR:CG	2.54	0.43
2:B:306:LEU:O	2:B:306:LEU:HD23	2.18	0.43
1:A:49:ASN:OD1	1:A:51:PHE:N	2.52	0.42
1:A:228:LEU:O	1:A:229:ASN:C	2.57	0.42
1:A:68:ASN:HA	1:A:69:PRO:HD2	1.83	0.42
1:A:-2:SER:HB2	1:A:-1:GLY:HA2	2.00	0.42
1:A:42:PHE:CZ	1:A:56:LEU:HD22	2.54	0.42
2:B:262:LEU:HD23	2:B:262:LEU:HA	1.74	0.42
1:A:12:PHE:O	1:A:15:ILE:N	2.53	0.42
1:A:190:ARG:O	1:A:191:YCM:C	2.67	0.42
1:A:7:MET:HE2	2:B:291:ARG:HH22	1.85	0.42
1:A:245:TYR:O	1:A:246:ILE:C	2.59	0.41
2:B:301:PHE:HA	2:B:304:MET:HE2	2.02	0.41
1:A:121:PHE:CE2	1:A:187:ARG:HA	2.56	0.41
1:A:12:PHE:HA	1:A:15:ILE:CG2	2.46	0.41
2:B:257:LEU:HD23	2:B:257:LEU:N	2.35	0.41
1:A:191:YCM:O	1:A:193:GLU:N	2.54	0.40
2:B:301:PHE:O	2:B:305:LEU:HB2	2.22	0.40
1:A:143:LEU:H	1:A:143:LEU:CD1	2.34	0.40
1:A:258:CYS:O	1:A:259:ILE:HD13	2.21	0.40
2:B:257:LEU:CD2	2:B:257:LEU:N	2.85	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	330/355 (93%)	277 (84%)	37 (11%)	16 (5%)	2	15
2	B	53/55 (96%)	42 (79%)	8 (15%)	3 (6%)	1	12
All	All	383/410 (93%)	319 (83%)	45 (12%)	19 (5%)	2	14

All (19) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	8	ASN
1	A	19	LEU
1	A	34	ALA
1	A	97	HIS
1	A	229	ASN
1	A	246	ILE
1	A	276	GLN
1	A	7	MET
1	A	33	LYS
1	A	37	VAL
1	A	136	THR
1	A	192	GLY
1	A	94	LYS
2	B	271	GLY
2	B	306	LEU
1	A	95	ASP
2	B	269	VAL
1	A	133	GLY
1	A	284	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	305/325 (94%)	259 (85%)	46 (15%)	3	12
2	B	48/48 (100%)	37 (77%)	11 (23%)	1	2

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
All	All	353/373 (95%)	296 (84%)	57 (16%)	2 10

All (57) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1	VAL
1	A	7	MET
1	A	9	ASN
1	A	18	MET
1	A	21	ILE
1	A	39	ASN
1	A	70	SER
1	A	75	VAL
1	A	77	GLU
1	A	79	LEU
1	A	91	GLU
1	A	94	LYS
1	A	95	ASP
1	A	113	LEU
1	A	114	ARG
1	A	130	ASN
1	A	134	GLN
1	A	138	GLU
1	A	143	LEU
1	A	146	GLN
1	A	162	ILE
1	A	173	ARG
1	A	176	ARG
1	A	183	LEU
1	A	193	GLU
1	A	197	LEU
1	A	207	LEU
1	A	216	GLU
1	A	223	TYR
1	A	229	ASN
1	A	233	HIS
1	A	240	SER
1	A	246	ILE
1	A	249	ILE
1	A	254	LYS
1	A	267	VAL
1	A	275	LEU

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Mol	Chain	Res	Type
1	A	279	LEU
1	A	286	GLU
1	A	290	LYS
1	A	292	LEU
1	A	298	LYS
1	A	315	ARG
1	A	318	GLN
1	A	320	GLN
1	A	329	GLN
2	B	255	ILE
2	B	257	LEU
2	B	263	LEU
2	B	276	LEU
2	B	282	ASN
2	B	291	ARG
2	B	293	HIS
2	B	295	MET
2	B	302	VAL
2	B	306	LEU
2	B	307	GLU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

1 non-standard protein/DNA/RNA residue 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
1	YCM	A	191	1,3	7,9,10	2.30	2 (28%)	4,10,12	1.90	1 (25%)

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
1	YCM	A	191	1,3	-	2/6/8/10	-

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	191	YCM	CE-NZ2	-4.76	1.17	1.32
1	A	191	YCM	OZ1-CE	3.03	1.33	1.24

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	191	YCM	CE-CD-SG	-3.32	103.83	113.59

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	A	191	YCM	SG-CD-CE-NZ2
1	A	191	YCM	SG-CD-CE-OZ1

There are no ring outliers.

1 monomer is involved in 7 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	A	191	YCM	7	0

5.5 Carbohydrates

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
3	NAG	C	1	1,3	14,14,15	0.85	1 (7%)	17,19,21	2.04	3 (17%)
3	NAG	C	2	3	14,14,15	1.16	2 (14%)	17,19,21	1.85	3 (17%)

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	NAG	C	1	1,3	-	3/6/23/26	0/1/1/1
3	NAG	C	2	3	-	4/6/23/26	0/1/1/1

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	C	2	NAG	O5-C1	-2.73	1.39	1.43
3	C	1	NAG	O5-C1	-2.73	1.39	1.43
3	C	2	NAG	O5-C5	-2.25	1.38	1.43

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	1	NAG	O5-C1-C2	-5.67	102.34	111.29
3	C	2	NAG	O5-C1-C2	-4.15	104.73	111.29
3	C	2	NAG	C2-N2-C7	-3.65	117.71	122.90
3	C	1	NAG	O5-C5-C6	3.53	112.74	107.20
3	C	2	NAG	C1-O5-C5	-3.20	107.86	112.19
3	C	1	NAG	C3-C4-C5	2.61	114.89	110.24

There are no chirality outliers.

All (7) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	C	2	NAG	C8-C7-N2-C2
3	C	2	NAG	O7-C7-N2-C2

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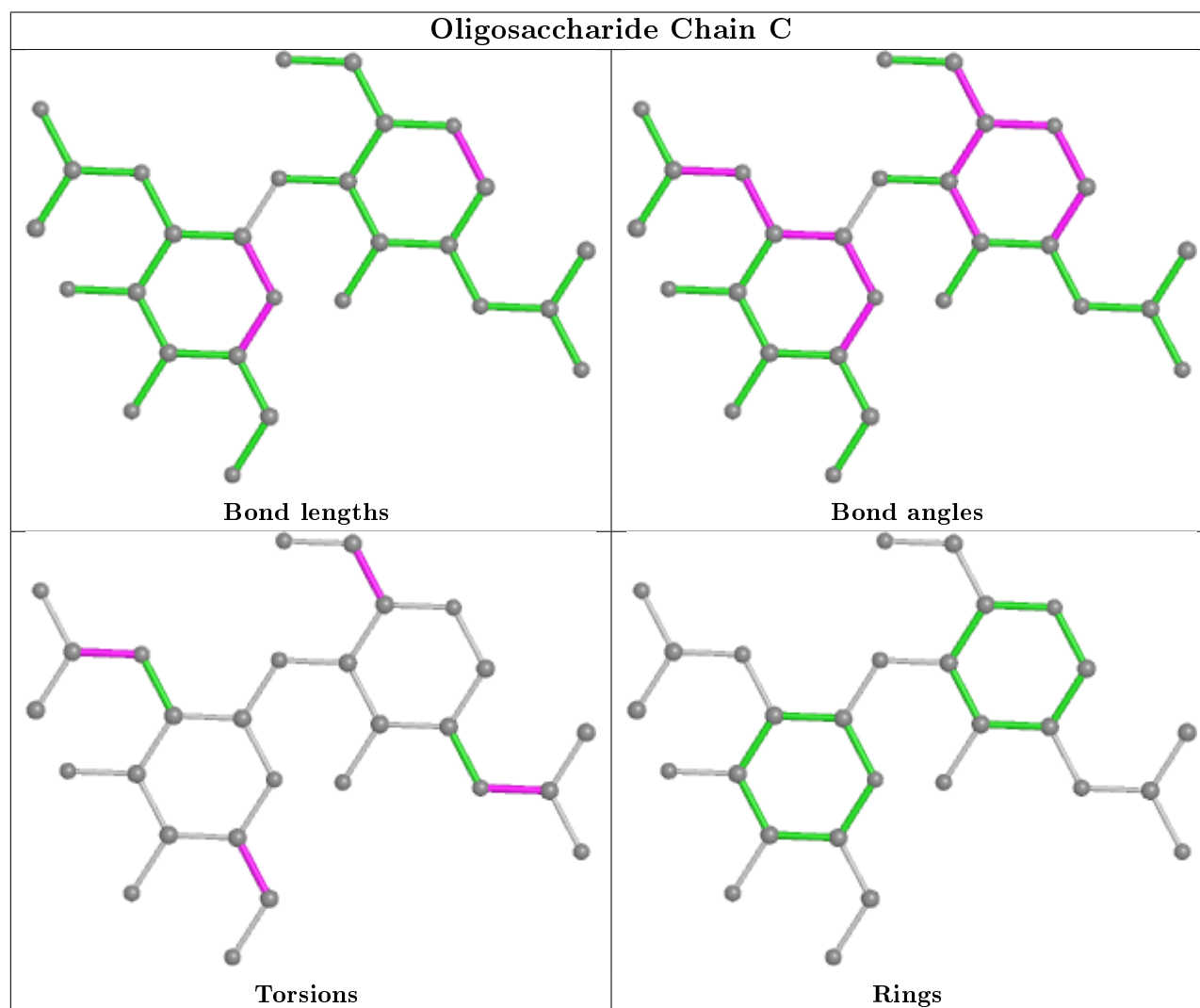
Mol	Chain	Res	Type	Atoms
3	C	1	NAG	C8-C7-N2-C2
3	C	1	NAG	O7-C7-N2-C2
3	C	2	NAG	C4-C5-C6-O6
3	C	2	NAG	O5-C5-C6-O6
3	C	1	NAG	O5-C5-C6-O6

There are no ring outliers.

2 monomers are involved in 10 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	C	2	NAG	4	0
3	C	1	NAG	9	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

Of 1 ligands modelled in this entry, 1 is 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	332/355 (93%)	-0.03	7 (2%) 63 62	36, 56, 94, 104	0
2	B	55/55 (100%)	0.14	2 (3%) 42 42	56, 73, 93, 95	0
All	All	387/410 (94%)	-0.01	9 (2%) 60 59	36, 59, 93, 104	0

All (9) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	98	ALA	4.2
1	A	5	SER	3.5
1	A	97	HIS	3.5
1	A	99	GLU	3.4
2	B	290	LEU	2.3
1	A	100	GLY	2.3
1	A	302	TYR	2.2
2	B	254	SER	2.2
1	A	4	GLY	2.1

6.2 Non-standard residues in protein, DNA, RNA chains [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.

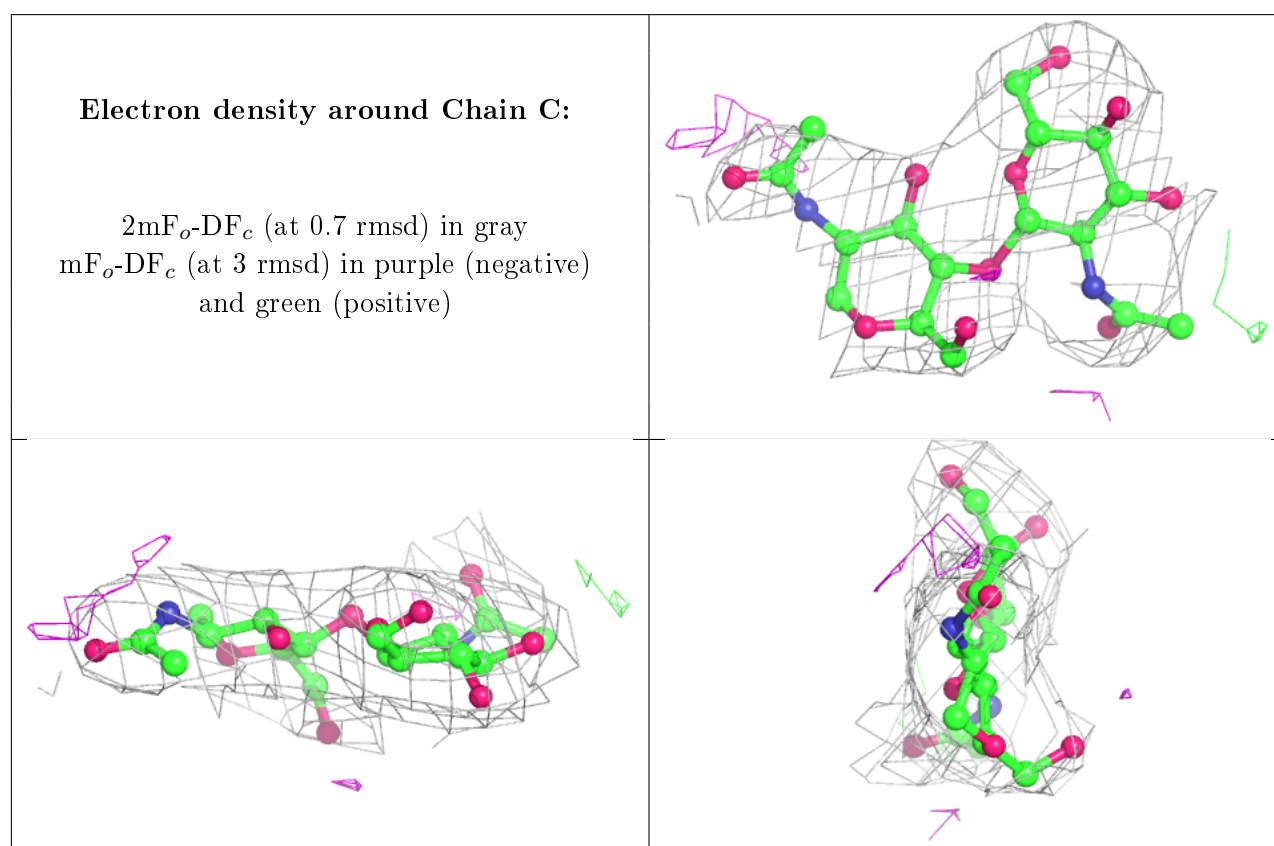
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
1	YCM	A	191	10/11	0.94	0.17	63,65,69,70	0

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
3	NAG	C	2	14/15	0.96	0.16	70,73,75,75	0
3	NAG	C	1	14/15	0.98	0.19	50,52,54,54	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(Å ²)	Q<0.9
4	ZN	A	344	1/1	1.00	0.12	77,77,77,77	0

6.5 Other polymers [i](#)

There are no such residues in this entry.