



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 12:24 AM UTC

PDB ID : 2V5M / pdb_00002v5m
Title : Structural basis for Dscam isoform specificity
Authors : Meijers, R.; Puettmann-Holgado, R.; Skiniotis, G.; Liu, J.-H.; Walz, T.;
Schmucker, D.; Wang, J.-H.
Deposited on : 2007-07-06
Resolution : 1.95 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

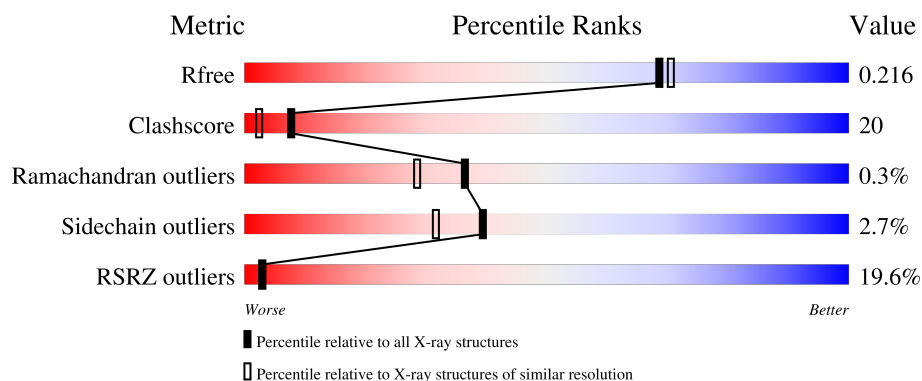
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.95 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	3494 (1.96-1.96)
Clashscore	190562	3612 (1.96-1.96)
Ramachandran outliers	187476	3587 (1.96-1.96)
Sidechain outliers	187428	3587 (1.96-1.96)
RSRZ outliers	180081	3495 (1.96-1.96)

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	388	20% 74% 23% .
2	B	2	50% 50%
2	C	2	50% 50%

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	GOL	A	601	-	X	-	-

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 3915 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 DSCAM.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	388	3144	1980	545	606	13	0	25	0

There are 20 discrepancies between the modelled and reference sequences:

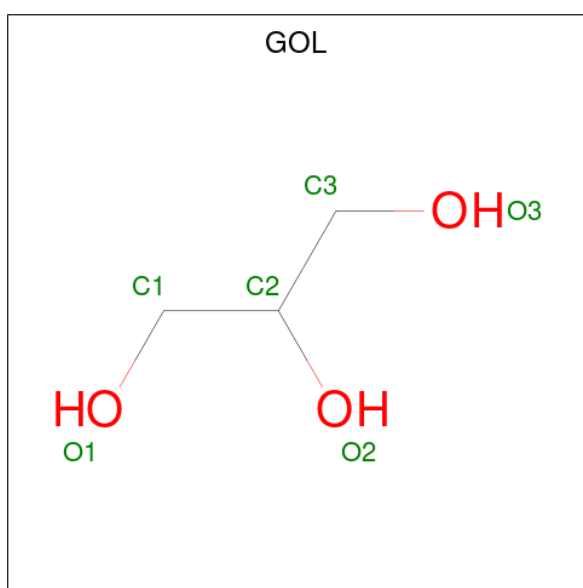
Chain	Residue	Modelled	Actual	Comment	Reference
A	207	ILE	VAL	variant	UNP Q9NBA1
A	211	ALA	PRO	variant	UNP Q9NBA1
A	213	ARG	LYS	variant	UNP Q9NBA1
A	214	THR	ILE	variant	UNP Q9NBA1
A	215	PRO	ASN	variant	UNP Q9NBA1
A	216	ALA	THR	variant	UNP Q9NBA1
A	218	VAL	THR	variant	UNP Q9NBA1
A	219	GLN	TYR	variant	UNP Q9NBA1
A	222	LEU	ASN	variant	UNP Q9NBA1
A	223	GLU	ILE	variant	UNP Q9NBA1
A	224	LEU	VAL	variant	UNP Q9NBA1
A	225	MET	GLU	variant	UNP Q9NBA1
A	226	VAL	SER	variant	UNP Q9NBA1
A	227	ALA	MET	variant	UNP Q9NBA1
A	228	HIS	ALA	variant	UNP Q9NBA1
A	229	THR	SER	variant	UNP Q9NBA1
A	230	ILE	THR	variant	UNP Q9NBA1
A	231	SER	ALA	variant	UNP Q9NBA1
A	232	LEU	ILE	variant	UNP Q9NBA1
A	239	PHE	TYR	variant	UNP Q9NBA1

- Molecule 2 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
2	B	2	Total	C	N	O	0	0	0
			28	16	2	10			
2	C	2	Total	C	N	O	0	0	0
			28	16	2	10			

- Molecule 3 is GLYCEROL (CCD ID: GOL) (formula: $C_3H_8O_3$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			6	3	3		

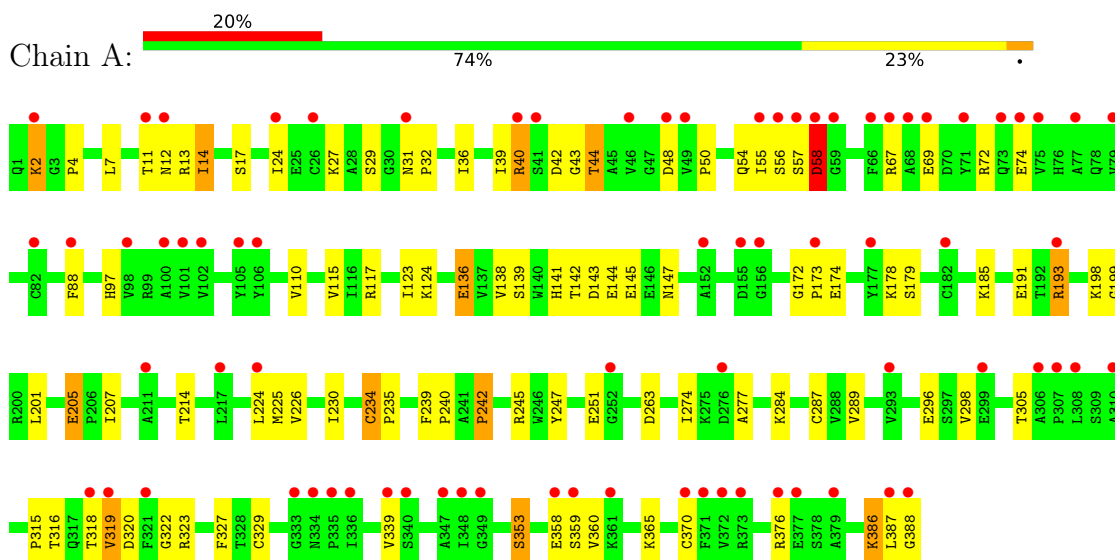
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	709	Total	O	0	0
			709	709		

3 Residue-property plots

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

• Molecule 1: DSCAM



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



• Molecule 2: 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 42 2 2	Depositor
Cell constants a, b, c, α , β , γ	99.17Å 99.17Å 163.99Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.00 – 1.95 20.00 – 1.95	Depositor EDS
% Data completeness (in resolution range)	98.6 (20.00-1.95) 98.6 (20.00-1.95)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.47 (at 1.95Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.169 , 0.205 0.180 , 0.216	Depositor DCC
R_{free} test set	3006 reflections (5.04%)	wwPDB-VP
Wilson B-factor (Å ²)	35.0	Xtriage
Anisotropy	0.377	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 52.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.35$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	3915	wwPDB-VP
Average B, all atoms (Å ²)	61.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.36% 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: GOL, NAG

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	0.78	4/3291 (0.1%)	1.38	27/4458 (0.6%)

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	1

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	386	LYS	C-O	8.54	1.34	1.23
1	A	387	LEU	C-O	7.98	1.33	1.23
1	A	386	LYS	C-N	7.50	1.44	1.33
1	A	320	ASP	C-O	6.53	1.31	1.23

All (27) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	193	ARG	NE-CZ-NH2	-11.24	109.08	119.20
1	A	193	ARG	NE-CZ-NH1	7.97	129.47	121.50
1	A	40	ARG	N-CA-CB	-7.84	98.23	110.06
1	A	88	PHE	N-CA-C	6.28	119.10	111.82
1	A	40	ARG	CA-CB-CG	6.07	126.24	114.10
1	A	242	PRO	N-CA-C	5.94	119.98	111.13
1	A	198	LYS	CD-CE-NZ	5.86	130.65	111.90
1	A	172	GLY	CA-C-N	5.85	125.99	119.32
1	A	172	GLY	C-N-CA	5.85	125.99	119.32

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	138	VAL	CA-CB-CG2	-5.55	100.97	110.40
1	A	17	SER	N-CA-CB	5.52	118.68	110.29
1	A	353	SER	N-CA-CB	-5.51	101.73	110.28
1	A	12	ASN	CB-CG-ND2	-5.44	108.24	116.40
1	A	14	ILE	N-CA-CB	-5.42	105.91	112.35
1	A	2	LYS	N-CA-C	5.35	117.22	108.55
1	A	2	LYS	CB-CA-C	-5.23	100.45	109.86
1	A	136[A]	GLU	CB-CG-CD	5.21	121.47	112.60
1	A	136[B]	GLU	CB-CG-CD	5.21	121.47	112.60
1	A	234	CYS	N-CA-CB	-5.15	104.70	111.09
1	A	386	LYS	CA-C-N	-5.12	113.31	122.07
1	A	386	LYS	C-N-CA	-5.12	113.31	122.07
1	A	205[A]	GLU	CA-C-N	-5.11	114.14	120.79
1	A	205[A]	GLU	C-N-CA	-5.11	114.14	120.79
1	A	205[B]	GLU	CA-C-N	-5.11	114.14	120.79
1	A	205[B]	GLU	C-N-CA	-5.11	114.14	120.79
1	A	72	ARG	NE-CZ-NH1	-5.08	116.42	121.50
1	A	72	ARG	NE-CZ-NH2	5.01	123.71	119.20

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	58[A]	ASP	Peptide

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	3144	0	3111	123	0
2	B	28	0	25	2	0
2	C	28	0	25	1	0
3	A	6	0	6	1	0
4	A	709	0	0	67	6
All	All	3915	0	3167	125	6

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 20.

All (125) 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:115:VAL:HB	4:A:2284:HOH:O	1.26	1.26
1:A:14:ILE:HG23	4:A:2037:HOH:O	1.50	1.10
1:A:57[B]:SER:O	1:A:58[B]:ASP:HB2	1.52	1.07
1:A:235:PRO:CD	4:A:2482:HOH:O	2.00	1.07
1:A:235:PRO:CG	4:A:2482:HOH:O	2.04	1.05
1:A:245[B]:ARG:HD2	4:A:2498:HOH:O	1.56	1.03
1:A:11:THR:OG1	4:A:2022:HOH:O	1.77	1.02
1:A:141:HIS:HD2	1:A:147:ASN:HD22	1.10	0.95
1:A:235:PRO:CB	4:A:2482:HOH:O	2.13	0.94
1:A:55:ILE:C	1:A:56[B]:SER:N	2.25	0.94
1:A:235:PRO:HD2	4:A:2482:HOH:O	1.61	0.92
4:A:2683:HOH:O	2:B:1:NAG:H81	1.73	0.89
1:A:191:GLU:HG3	4:A:2152:HOH:O	1.73	0.88
1:A:141:HIS:HD2	1:A:147:ASN:ND2	1.73	0.86
1:A:57[B]:SER:O	1:A:58[B]:ASP:CB	2.23	0.85
1:A:245[B]:ARG:HG2	1:A:247:TYR:CE2	2.12	0.84
1:A:67:ARG:HA	2:B:1:NAG:H83	1.60	0.83
1:A:42:ASP:OD2	1:A:44:THR:HG22	1.79	0.83
1:A:274:ILE:CG2	4:A:2554:HOH:O	2.30	0.80
1:A:225[B]:MET:HE2	1:A:305:THR:OG1	1.82	0.80
1:A:57[A]:SER:O	1:A:58[A]:ASP:CB	2.30	0.80
1:A:245[B]:ARG:HG2	1:A:247:TYR:CZ	2.17	0.79
1:A:144[B]:GLU:OE1	4:A:2347:HOH:O	1.99	0.79
1:A:225[B]:MET:CE	1:A:305:THR:OG1	2.32	0.77
1:A:141:HIS:CD2	1:A:147:ASN:HD22	1.98	0.76
1:A:388:GLY:C	4:A:2598:HOH:O	2.28	0.76
1:A:144[B]:GLU:HB2	1:A:178[B]:LYS:HE3	1.67	0.76
1:A:358[B]:GLU:OE2	4:A:2654:HOH:O	2.05	0.75
1:A:235:PRO:HB2	4:A:2482:HOH:O	1.80	0.74
1:A:145[B]:GLU:HG3	4:A:2350:HOH:O	1.86	0.74
1:A:245[B]:ARG:CD	4:A:2498:HOH:O	2.25	0.74
1:A:319:VAL:HG22	1:A:360:VAL:HG21	1.69	0.73
1:A:315:PRO:HD2	4:A:2592:HOH:O	1.87	0.73
1:A:329:CYS:HG	1:A:370:CYS:HG	1.33	0.73
1:A:319:VAL:CG2	1:A:360:VAL:HG21	2.20	0.72
1:A:14:ILE:HD11	4:A:2022:HOH:O	1.90	0.71
1:A:42:ASP:OD2	1:A:42:ASP:C	2.34	0.71
1:A:245[B]:ARG:CG	1:A:247:TYR:CE2	2.75	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:13[A]:ARG:NH1	4:A:2028:HOH:O	2.21	0.70
1:A:376:ARG:HD2	4:A:2030:HOH:O	1.92	0.69
1:A:57[A]:SER:O	1:A:58[A]:ASP:HB2	1.93	0.69
1:A:296:GLU:OE2	4:A:2567:HOH:O	2.10	0.68
1:A:263[B]:ASP:OD1	4:A:2527:HOH:O	2.12	0.68
1:A:143[A]:ASP:OD2	1:A:179:SER:HB3	1.95	0.66
1:A:54:GLN:HG2	4:A:2115:HOH:O	1.94	0.65
1:A:143[B]:ASP:CG	4:A:2341:HOH:O	2.38	0.65
1:A:274:ILE:HG21	4:A:2554:HOH:O	1.93	0.65
1:A:14:ILE:CG2	4:A:2037:HOH:O	2.24	0.63
1:A:13[A]:ARG:HD3	4:A:2028:HOH:O	1.99	0.62
1:A:97:HIS:HD2	4:A:2010:HOH:O	1.82	0.61
1:A:143[B]:ASP:OD1	1:A:179:SER:N	2.29	0.60
1:A:230[B]:ILE:HD12	4:A:2554:HOH:O	2.00	0.60
1:A:274:ILE:HG22	4:A:2554:HOH:O	1.99	0.59
1:A:318:THR:HA	1:A:386:LYS:O	2.02	0.59
1:A:50:PRO:HG3	4:A:2120:HOH:O	2.03	0.58
1:A:11:THR:HG22	4:A:2008:HOH:O	2.04	0.58
1:A:13[B]:ARG:NH2	4:A:2030:HOH:O	2.36	0.58
1:A:11:THR:N	4:A:2022:HOH:O	2.36	0.57
1:A:144[B]:GLU:HB2	1:A:178[B]:LYS:CE	2.36	0.56
1:A:245[B]:ARG:HG2	1:A:247:TYR:OH	2.05	0.56
1:A:14:ILE:HA	4:A:2037:HOH:O	2.06	0.55
1:A:245[B]:ARG:CG	1:A:247:TYR:CZ	2.89	0.55
1:A:323:ARG:NH1	4:A:2601:HOH:O	2.37	0.54
1:A:178[B]:LYS:HG3	4:A:2341:HOH:O	2.08	0.54
1:A:298:VAL:HG11	4:A:2480:HOH:O	2.07	0.54
1:A:42:ASP:OD2	1:A:43:GLY:N	2.40	0.54
1:A:234:CYS:HG	1:A:287:CYS:CB	2.20	0.54
1:A:365:LYS:HE2	4:A:2659:HOH:O	2.08	0.53
1:A:67:ARG:HD3	4:A:2162:HOH:O	2.06	0.53
1:A:117[B]:ARG:HG2	1:A:173:PRO:HD3	1.91	0.52
1:A:224:LEU:HB3	4:A:2472:HOH:O	2.08	0.52
1:A:245[B]:ARG:HD3	1:A:247:TYR:CZ	2.44	0.52
1:A:74:GLU:HG2	4:A:2186:HOH:O	2.11	0.50
1:A:316:THR:CG2	4:A:2004:HOH:O	2.58	0.50
1:A:143[B]:ASP:OD1	4:A:2341:HOH:O	2.19	0.50
1:A:44:THR:HG21	4:A:2093:HOH:O	2.11	0.49
1:A:319:VAL:CG2	1:A:360:VAL:CG2	2.90	0.49
1:A:57[A]:SER:O	1:A:58[A]:ASP:HB3	2.09	0.49
1:A:230[A]:ILE:HD11	4:A:2472:HOH:O	2.12	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:13[B]:ARG:CZ	4:A:2030:HOH:O	2.60	0.48
1:A:242:PRO:HB2	1:A:289[A]:VAL:HG23	1.94	0.48
1:A:284:LYS:CE	4:A:2556:HOH:O	2.61	0.48
1:A:56[B]:SER:HB3	4:A:2135:HOH:O	2.13	0.48
1:A:7:LEU:HD21	4:A:2069:HOH:O	2.14	0.47
1:A:277:ALA:HB2	4:A:2554:HOH:O	2.14	0.47
4:A:2708:HOH:O	2:C:2:NAG:O4	2.08	0.47
1:A:322:GLY:HA2	1:A:359:SER:OG	2.15	0.47
1:A:365:LYS:HG3	4:A:2659:HOH:O	2.14	0.46
1:A:234:CYS:O	3:A:601:GOL:H32	2.16	0.46
1:A:13[A]:ARG:NH1	4:A:2029:HOH:O	2.49	0.46
1:A:69[B]:GLU:CD	4:A:2162:HOH:O	2.58	0.46
1:A:4:PRO:HA	1:A:29:SER:O	2.15	0.46
1:A:2:LYS:NZ	4:A:2004:HOH:O	2.45	0.45
1:A:54:GLN:HG2	4:A:2119:HOH:O	2.17	0.45
1:A:339:VAL:HG13	1:A:370:CYS:SG	2.56	0.45
1:A:358[B]:GLU:N	1:A:358[B]:GLU:CD	2.75	0.45
1:A:298:VAL:CG1	4:A:2480:HOH:O	2.66	0.44
1:A:31:ASN:HD22	1:A:32:PRO:HA	1.82	0.44
1:A:144[B]:GLU:CB	1:A:178[B]:LYS:HE3	2.43	0.44
1:A:277:ALA:CB	4:A:2554:HOH:O	2.65	0.44
1:A:123:ILE:HD11	1:A:201:LEU:HD21	1.99	0.44
1:A:388:GLY:C	4:A:2682:HOH:O	2.60	0.43
1:A:39:ILE:HD12	1:A:43:GLY:HA2	2.01	0.42
1:A:225[B]:MET:HG3	1:A:226:VAL:N	2.34	0.42
1:A:24:ILE:HD13	1:A:24:ILE:HG21	1.74	0.42
1:A:207:ILE:HD12	4:A:2163:HOH:O	2.19	0.42
1:A:245[B]:ARG:HG3	1:A:247:TYR:CE2	2.54	0.42
1:A:327:PHE:HE1	4:A:2593:HOH:O	2.01	0.42
1:A:142:THR:HB	1:A:178[B]:LYS:HE2	2.00	0.42
1:A:358[B]:GLU:CD	1:A:358[B]:GLU:H	2.28	0.42
1:A:139:SER:OG	1:A:141:HIS:HE1	2.02	0.41
1:A:40:ARG:CD	4:A:2100:HOH:O	2.68	0.41
1:A:142:THR:CB	1:A:178[B]:LYS:HE2	2.51	0.41
1:A:205[A]:GLU:OE2	4:A:2438:HOH:O	2.22	0.41
1:A:214:THR:HB	4:A:2453:HOH:O	2.19	0.41
1:A:239:PHE:HA	1:A:240:PRO:C	2.46	0.41
1:A:40:ARG:HD3	4:A:2100:HOH:O	2.20	0.41
1:A:242:PRO:HB2	1:A:289[A]:VAL:CG2	2.50	0.40
1:A:136[B]:GLU:HG2	1:A:185:LYS:O	2.21	0.40
1:A:173:PRO:HD2	1:A:174:GLU:OE1	2.22	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:55:ILE:O	1:A:56[B]:SER:N	2.52	0.40
1:A:110:VAL:HG12	1:A:199:GLY:HA3	2.02	0.40

All (6) 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:2185:HOH:O	4:A:2211:HOH:O[8_665]	1.50	0.70
4:A:2400:HOH:O	4:A:2400:HOH:O[2_765]	1.74	0.46
4:A:2029:HOH:O	4:A:2450:HOH:O[5_755]	1.82	0.38
4:A:2429:HOH:O	4:A:2429:HOH:O[6_565]	2.00	0.20
4:A:2279:HOH:O	4:A:2316:HOH:O[5_755]	2.02	0.18
4:A:2198:HOH:O	4:A:2267:HOH:O[6_555]	2.04	0.16

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	411/388 (106%)	398 (97%)	11 (3%)	2 (0%)	24 16

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	58[A]	ASP
1	A	58[B]	ASP

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	356/332 (107%)	347 (98%)	9 (2%)	42 34

All (9) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	27	LYS
1	A	36	ILE
1	A	44	THR
1	A	48	ASP
1	A	124	LYS
1	A	193	ARG
1	A	251	GLU
1	A	319	VAL
1	A	353	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	12	ASN
1	A	31	ASN
1	A	97	HIS
1	A	111	ASN
1	A	141	HIS
1	A	147	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](#)

4 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	NAG	B	1	1,2	14,14,15	0.86	0	17,19,21	3.12	9 (52%)
2	NAG	B	2	2	14,14,15	0.58	0	17,19,21	2.27	8 (47%)
2	NAG	C	1	1,2	14,14,15	0.41	0	17,19,21	2.67	7 (41%)
2	NAG	C	2	2	14,14,15	0.49	0	17,19,21	1.28	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
2	NAG	B	1	1,2	-	4/6/23/26	0/1/1/1
2	NAG	B	2	2	-	2/6/23/26	0/1/1/1
2	NAG	C	1	1,2	-	2/6/23/26	0/1/1/1
2	NAG	C	2	2	-	0/6/23/26	0/1/1/1

There are no bond length outliers.

All (27) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	1	NAG	O6-C6-C5	-6.91	87.82	111.33
2	C	1	NAG	C1-O5-C5	5.86	120.03	112.19
2	C	1	NAG	C1-C2-N2	5.53	119.15	110.43
2	B	1	NAG	C3-C4-C5	5.46	120.12	110.23
2	C	1	NAG	C6-C5-C4	-5.01	100.72	113.02
2	B	1	NAG	C6-C5-C4	-4.28	102.51	113.02
2	B	1	NAG	C1-O5-C5	4.10	117.68	112.19
2	B	2	NAG	C2-N2-C7	-3.98	117.57	122.90
2	B	1	NAG	C4-C3-C2	-3.56	105.81	111.02
2	B	2	NAG	C4-C3-C2	-3.39	106.05	111.02
2	B	2	NAG	C6-C5-C4	-3.37	104.75	113.02
2	B	1	NAG	C2-N2-C7	3.26	127.26	122.90
2	B	2	NAG	C1-O5-C5	-3.15	107.96	112.19

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	1	NAG	O7-C7-C8	-3.14	116.45	122.05
2	B	2	NAG	C1-C2-N2	3.10	115.32	110.43
2	B	2	NAG	O5-C5-C6	2.95	113.41	107.66
2	C	1	NAG	O5-C5-C6	-2.69	102.43	107.66
2	B	1	NAG	O7-C7-N2	2.56	126.51	121.98
2	C	2	NAG	C1-C2-N2	2.37	114.17	110.43
2	C	1	NAG	O5-C1-C2	-2.33	107.69	111.29
2	B	2	NAG	O4-C4-C5	2.29	114.97	109.32
2	C	1	NAG	O4-C4-C3	-2.29	104.99	110.38
2	C	2	NAG	C4-C3-C2	-2.15	107.87	111.02
2	C	1	NAG	C3-C4-C5	2.15	114.13	110.23
2	B	2	NAG	O7-C7-N2	2.15	125.78	121.98
2	B	1	NAG	O5-C1-C2	-2.14	107.98	111.29
2	C	2	NAG	O7-C7-N2	2.04	125.58	121.98

There are no chirality outliers.

All (8) torsion outliers are listed below:

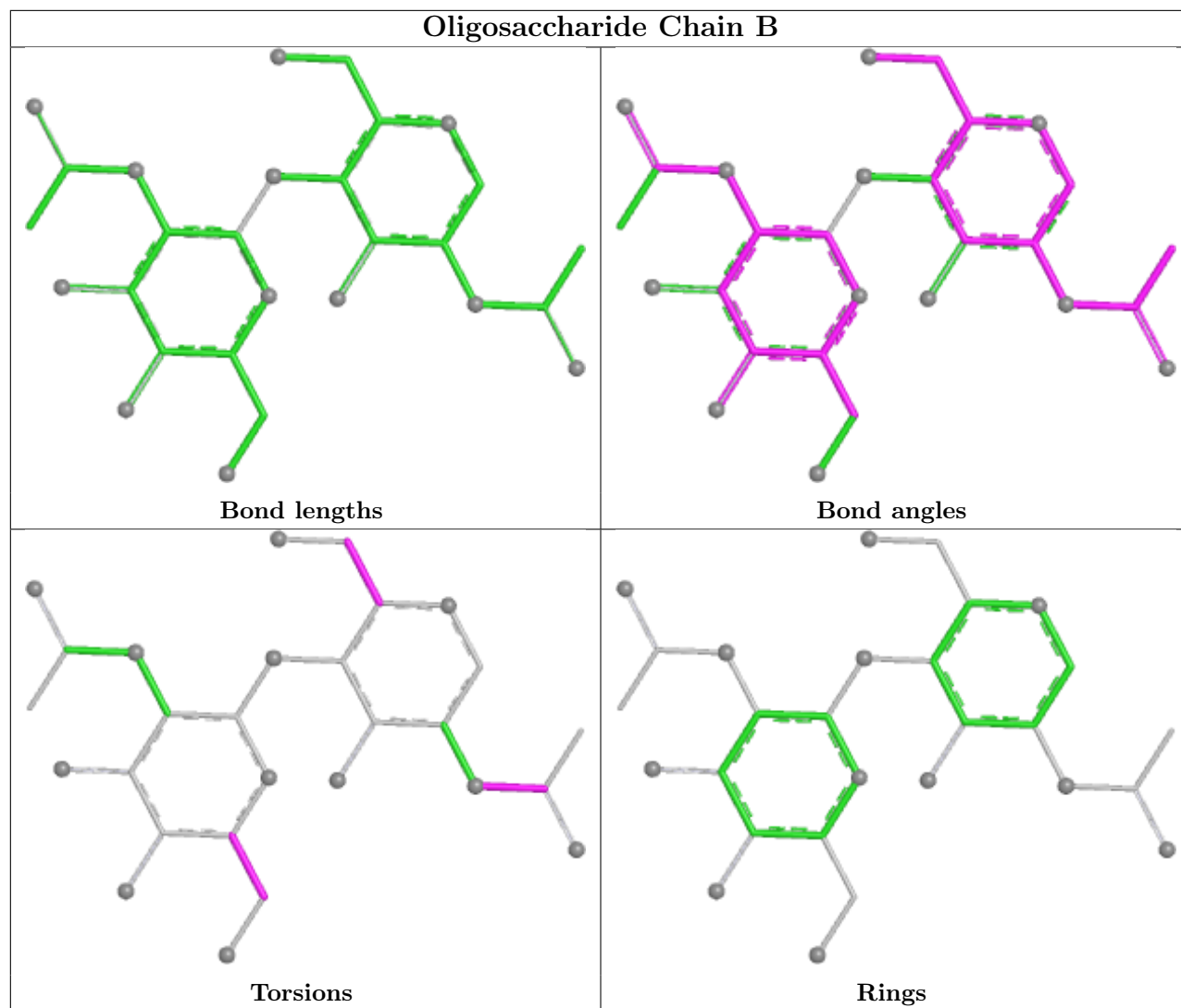
Mol	Chain	Res	Type	Atoms
2	C	1	NAG	O5-C5-C6-O6
2	B	1	NAG	C4-C5-C6-O6
2	C	1	NAG	C4-C5-C6-O6
2	B	1	NAG	O5-C5-C6-O6
2	B	1	NAG	C8-C7-N2-C2
2	B	2	NAG	C4-C5-C6-O6
2	B	2	NAG	O5-C5-C6-O6
2	B	1	NAG	O7-C7-N2-C2

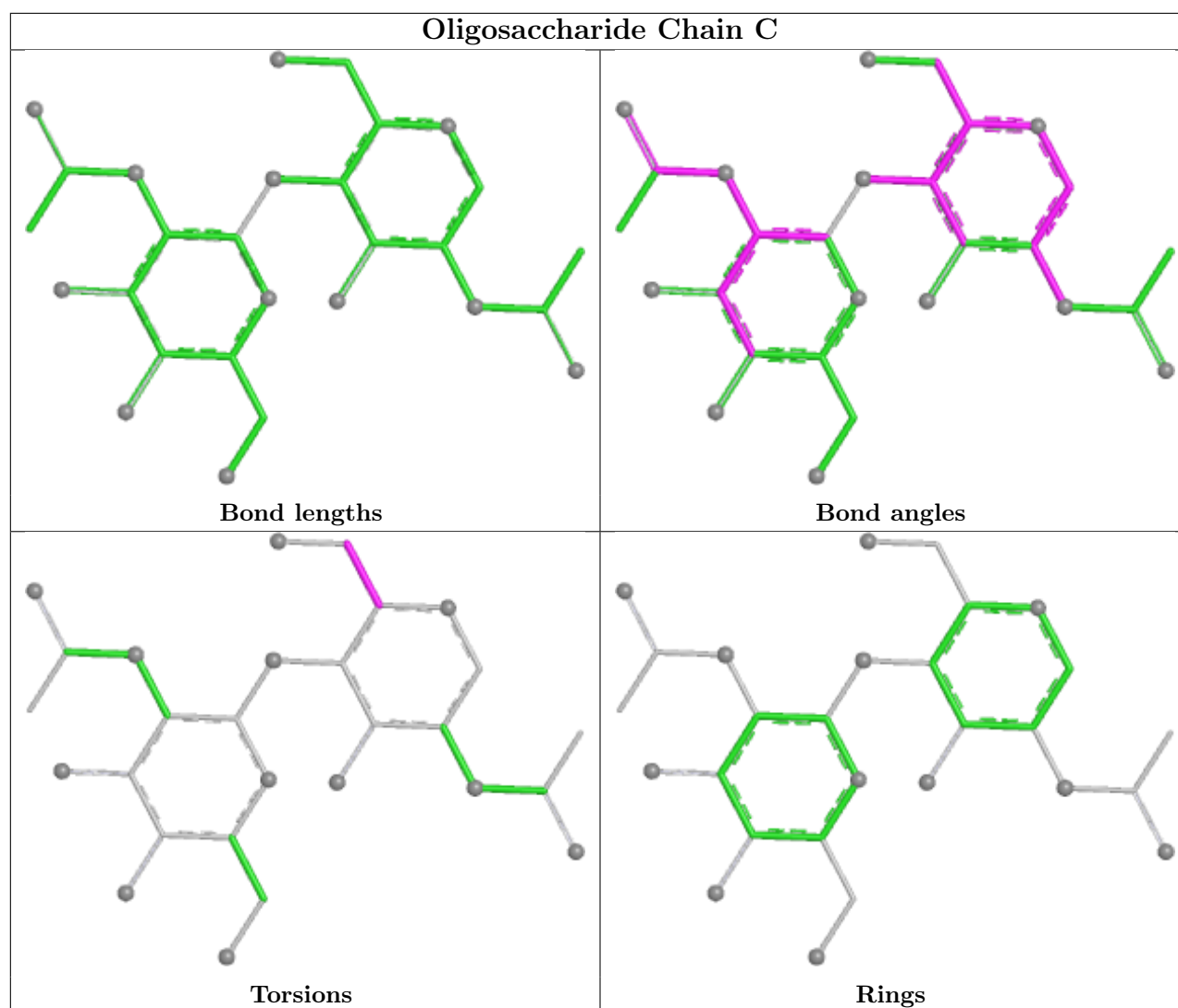
There are no ring outliers.

2 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	1	NAG	2	0
2	C	2	NAG	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](#)

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	GOL	A	601	-	5,5,5	0.50	0	5,5,5	5.98	3 (60%)

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	GOL	A	601	-	-	4/4/4/4	-

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	601	GOL	C3-C2-C1	-12.25	66.86	111.80
3	A	601	GOL	O2-C2-C1	4.50	127.83	109.18
3	A	601	GOL	O2-C2-C3	2.27	118.58	109.18

There are no chirality outliers.

All (4) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	601	GOL	O1-C1-C2-C3
3	A	601	GOL	C1-C2-C3-O3
3	A	601	GOL	O1-C1-C2-O2
3	A	601	GOL	O2-C2-C3-O3

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	601	GOL	1	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

The following chains have linkage breaks:

Mol	Chain	Number of breaks
1	A	1

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	A	55:ILE	C	56[B]:SER	N	2.25

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, 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	388/388 (100%)	1.45	76 (19%) 3 3	32, 57, 68, 81	25 (6%)

All (76) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	59[A]	GLY	6.1
1	A	58[A]	ASP	5.3
1	A	41	SER	4.3
1	A	177	TYR	4.2
1	A	359	SER	3.8
1	A	55	ILE	3.7
1	A	370	CYS	3.5
1	A	319	VAL	3.5
1	A	173	PRO	3.4
1	A	182	CYS	3.4
1	A	376	ARG	3.3
1	A	321	PHE	3.1
1	A	388	GLY	3.0
1	A	88	PHE	2.8
1	A	371	PHE	2.8
1	A	349	GLY	2.8
1	A	101	VAL	2.8
1	A	293	VAL	2.8
1	A	57[A]	SER	2.8
1	A	82	CYS	2.8
1	A	69[A]	GLU	2.8
1	A	306	ALA	2.7
1	A	252	GLY	2.7
1	A	372	VAL	2.6
1	A	155	ASP	2.6
1	A	79	VAL	2.6
1	A	26	CYS	2.6

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Mol	Chain	Res	Type	RSRZ
1	A	308	LEU	2.6
1	A	276	ASP	2.5
1	A	334	ASN	2.5
1	A	339	VAL	2.5
1	A	348	ILE	2.5
1	A	11	THR	2.5
1	A	361	LYS	2.5
1	A	74	GLU	2.4
1	A	373	ARG	2.4
1	A	49	VAL	2.4
1	A	40	ARG	2.3
1	A	152	ALA	2.3
1	A	333	GLY	2.3
1	A	211	ALA	2.3
1	A	379	ALA	2.3
1	A	56[A]	SER	2.3
1	A	67	ARG	2.3
1	A	48	ASP	2.3
1	A	340	SER	2.3
1	A	73	GLN	2.3
1	A	66	PHE	2.2
1	A	318	THR	2.2
1	A	217	LEU	2.2
1	A	299	GLU	2.2
1	A	387	LEU	2.2
1	A	307	PRO	2.2
1	A	71	TYR	2.2
1	A	106	TYR	2.2
1	A	24	ILE	2.2
1	A	98	VAL	2.2
1	A	358[A]	GLU	2.2
1	A	100	ALA	2.2
1	A	336	ILE	2.2
1	A	377	GLU	2.2
1	A	335	PRO	2.2
1	A	2	LYS	2.2
1	A	46	VAL	2.1
1	A	75	VAL	2.1
1	A	156	GLY	2.1
1	A	224	LEU	2.1
1	A	310	ALA	2.1
1	A	105	TYR	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	31	ASN	2.1
1	A	12	ASN	2.1
1	A	68	ALA	2.1
1	A	77	ALA	2.1
1	A	102	VAL	2.0
1	A	347	ALA	2.0
1	A	193	ARG	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](#)

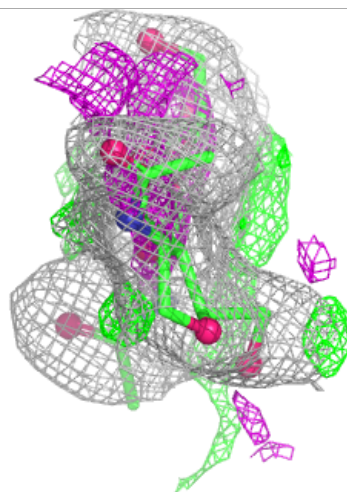
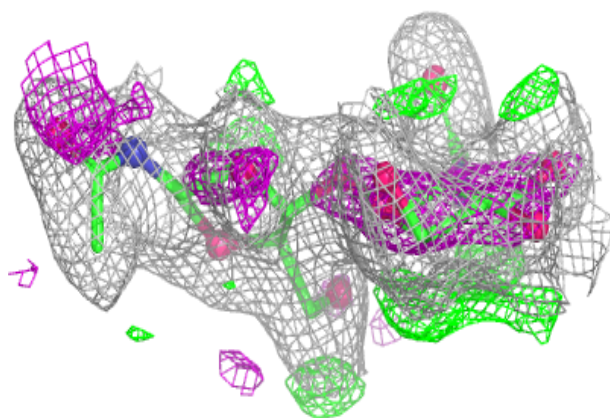
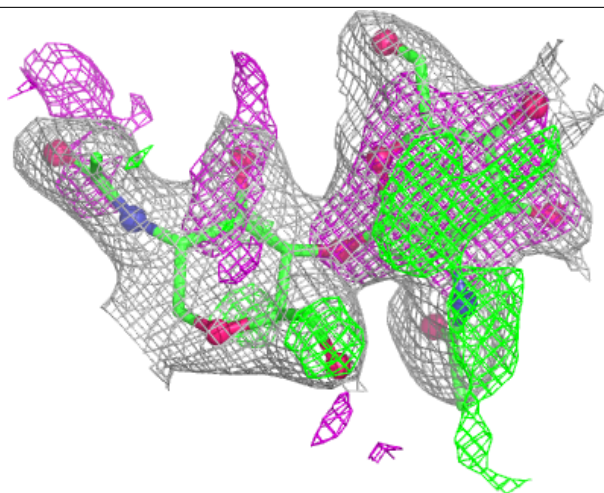
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

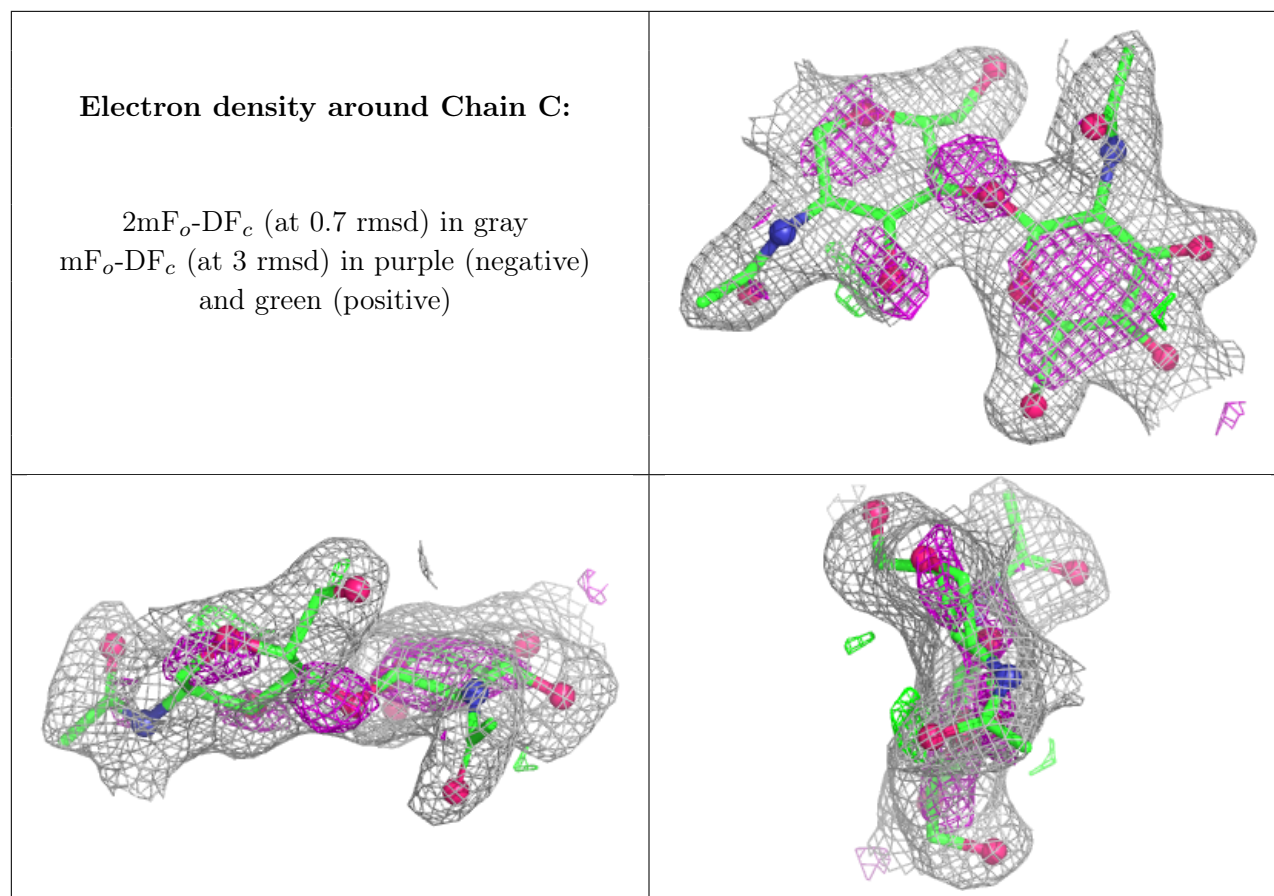
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	NAG	B	2	14/15	0.71	0.19	59,66,74,80	0
2	NAG	C	1	14/15	0.84	0.17	62,69,77,84	0
2	NAG	C	2	14/15	0.84	0.16	64,68,83,85	0
2	NAG	B	1	14/15	0.88	0.14	43,53,62,71	0

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.

Electron density around Chain B:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)





6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	GOL	A	601	6/6	0.92	0.24	40,46,53,53	0

6.5 Other polymers [i](#)

There are no such residues in this entry.