



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 1, 2026 – 11:52 PM UTC

PDB ID : 1Q25 / pdb_00001q25
Title : Crystal structure of N-terminal 3 domains of CI-MPR
Authors : Olson, L.J.; Dahms, N.M.; Kim, J.-J.P.
Deposited on : 2003-07-23
Resolution : 1.80 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

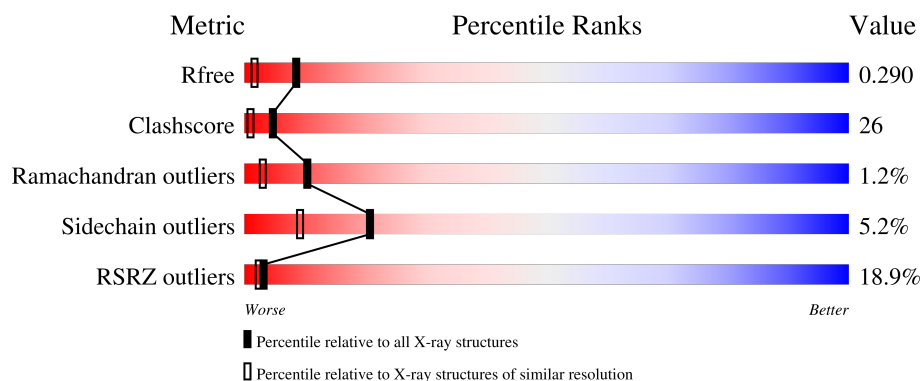
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.80 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	7662 (1.80-1.80)
Clashscore	190562	8479 (1.80-1.80)
Ramachandran outliers	187476	8391 (1.80-1.80)
Sidechain outliers	187428	8390 (1.80-1.80)
RSRZ outliers	180081	7663 (1.80-1.80)

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	432	19% 66% 28% 5%
2	B	5	80% 20%
3	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
4	GOL	A	4005	-	X	-	-

2 Entry composition [i](#)

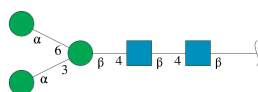
There are 5 unique types of molecules in this entry. The entry contains 3763 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 cation-independent mannose 6-phosphate receptor.

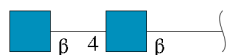
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	428	Total	C	N	O	S	0	0	0
			3303	2060	564	651	28			

- Molecule 2 is an oligosaccharide called alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-6)]beta-D-mannopyranose-(1-4)-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	5	Total	C	N	O	0	0	0
			61	34	2	25			

- 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 GLYCEROL (CCD ID: GOL) (formula: C₃H₈O₃).



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

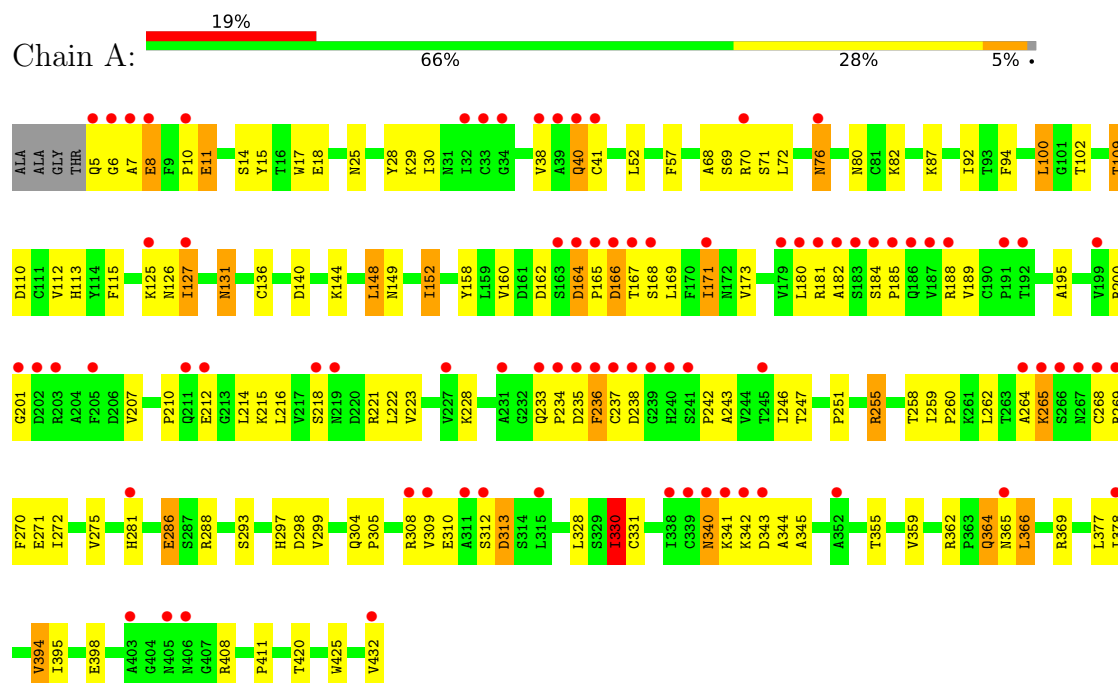
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	365	Total	O	0	0
			365	365		

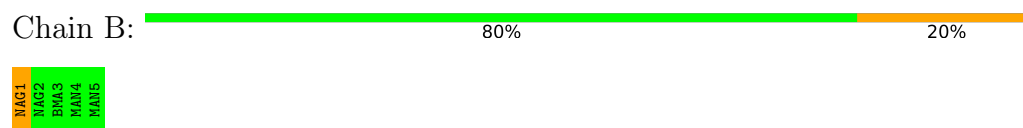
3 Residue-property plots

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: cation-independent mannose 6-phosphate receptor



- Molecule 2: alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-6)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose



- 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 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	61.16Å 84.75Å 96.05Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	19.94 – 1.80 19.94 – 1.80	Depositor EDS
% Data completeness (in resolution range)	87.1 (19.94-1.80) 87.0 (19.94-1.80)	Depositor EDS
R_{merge}	0.04	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.03 (at 1.76Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.235 , 0.283 0.245 , 0.290	Depositor DCC
R_{free} test set	2820 reflections (6.72%)	wwPDB-VP
Wilson B-factor (Å ²)	21.9	Xtriage
Anisotropy	0.257	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 47.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	3763	wwPDB-VP
Average B, all atoms (Å ²)	33.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.19% 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: NAG, MAN, GOL, BMA

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	0.39	0/3371	0.97	13/4570 (0.3%)

There are no bond length outliers.

All (13) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	76	ASN	CB-CG-ND2	7.85	128.17	116.40
1	A	76	ASN	CA-CB-CG	7.37	119.97	112.60
1	A	166	ASP	N-CA-C	-7.24	105.37	114.56
1	A	345	ALA	N-CA-C	-7.01	103.71	112.90
1	A	76	ASN	CB-CA-C	-6.47	97.54	111.78
1	A	286	GLU	N-CA-C	6.38	119.42	109.52
1	A	109	THR	N-CA-C	-6.02	100.48	109.94
1	A	236	PHE	N-CA-C	-6.02	105.78	113.01
1	A	330	ILE	N-CA-C	5.47	117.45	108.97
1	A	131	ASN	N-CA-C	-5.29	105.51	111.28
1	A	355	THR	N-CA-C	-5.24	106.89	113.28
1	A	420	THR	N-CA-C	-5.17	100.09	108.52
1	A	112	VAL	N-CA-C	5.00	115.69	108.48

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	3303	0	3159	171	0
2	B	61	0	52	1	0
3	C	28	0	25	2	0
4	A	6	0	4	0	0
5	A	365	0	0	11	0
All	All	3763	0	3240	171	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 26.

All (171) 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:92:ILE:HD13	1:A:115:PHE:HB2	1.37	1.04
1:A:246:ILE:HD13	1:A:272:ILE:HB	1.49	0.94
1:A:309:VAL:HG11	1:A:313:ASP:OD1	1.71	0.91
1:A:228:LYS:HB3	1:A:228:LYS:HZ2	1.33	0.90
1:A:164:ASP:HB3	1:A:167:THR:OG1	1.73	0.88
1:A:378:ILE:CD1	1:A:394:VAL:HB	2.04	0.88
1:A:80:ASN:HD21	1:A:87:LYS:HZ3	1.22	0.87
1:A:109:THR:HG22	1:A:110:ASP:H	1.43	0.84
1:A:195:ALA:HB2	1:A:210:PRO:HD3	1.57	0.84
1:A:293:SER:H	1:A:297:HIS:HD2	1.24	0.84
1:A:149:ASN:HA	1:A:152:ILE:HD13	1.59	0.84
1:A:309:VAL:HG13	1:A:408:ARG:NH2	1.94	0.81
1:A:76:ASN:OD1	5:A:774:HOH:O	1.99	0.79
1:A:80:ASN:HD21	1:A:87:LYS:NZ	1.81	0.77
1:A:207:VAL:HG23	1:A:242:PRO:HB3	1.67	0.76
1:A:299:VAL:HG23	1:A:432:VAL:HG13	1.68	0.75
1:A:92:ILE:HD13	1:A:115:PHE:CB	2.16	0.74
1:A:160:VAL:CG2	1:A:171:ILE:HD12	2.18	0.73
1:A:340:ASN:CG	1:A:341:LYS:H	1.97	0.73
1:A:228:LYS:HB3	1:A:228:LYS:NZ	2.03	0.73
1:A:185:PRO:HA	1:A:188:ARG:HE	1.53	0.72
1:A:309:VAL:CG1	1:A:313:ASP:OD1	2.36	0.72
1:A:236:PHE:CE1	1:A:237:CYS:SG	2.82	0.72
1:A:6:GLY:HA2	1:A:10:PRO:CG	2.21	0.70
1:A:80:ASN:ND2	1:A:87:LYS:NZ	2.40	0.70
1:A:14:SER:HB2	1:A:15:TYR:CE1	2.27	0.69
1:A:330:ILE:HG23	1:A:331:CYS:SG	2.31	0.69
1:A:109:THR:HG22	1:A:110:ASP:N	2.07	0.69
1:A:40:GLN:NE2	1:A:40:GLN:H	1.92	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:242:PRO:HA	1:A:268:CYS:O	1.93	0.68
1:A:184:SER:O	1:A:188:ARG:HG3	1.94	0.68
1:A:246:ILE:CD1	1:A:272:ILE:HB	2.21	0.68
1:A:40:GLN:H	1:A:40:GLN:HE21	1.40	0.68
1:A:215:LYS:HE2	1:A:223:VAL:HG23	1.76	0.68
1:A:411:PRO:HG3	1:A:425:TRP:CZ3	2.29	0.67
1:A:149:ASN:CA	1:A:152:ILE:HD13	2.25	0.66
1:A:25:ASN:ND2	1:A:52:LEU:HB2	2.12	0.65
1:A:341:LYS:HD3	1:A:341:LYS:C	2.21	0.65
1:A:364:GLN:NE2	1:A:364:GLN:H	1.94	0.65
1:A:364:GLN:H	1:A:364:GLN:HE21	1.44	0.65
1:A:125:LYS:C	1:A:127:ILE:H	2.04	0.64
1:A:233:GLN:HG3	1:A:242:PRO:HD3	1.79	0.64
1:A:160:VAL:HG23	1:A:171:ILE:HD12	1.77	0.64
1:A:251:PRO:HG3	1:A:275:VAL:CG1	2.28	0.63
1:A:136:CYS:SG	1:A:152:ILE:HD11	2.39	0.62
1:A:378:ILE:HD13	1:A:394:VAL:HB	1.80	0.61
1:A:242:PRO:HG3	1:A:268:CYS:SG	2.40	0.61
1:A:242:PRO:HA	1:A:268:CYS:HB3	1.84	0.60
1:A:233:GLN:HG3	1:A:242:PRO:CD	2.32	0.59
1:A:411:PRO:HG3	1:A:425:TRP:CE3	2.36	0.59
1:A:207:VAL:HG22	1:A:207:VAL:O	2.03	0.59
1:A:11:GLU:HG3	5:A:697:HOH:O	2.02	0.58
1:A:340:ASN:ND2	1:A:341:LYS:H	2.00	0.58
1:A:71:SER:HB2	1:A:94:PHE:O	2.04	0.58
1:A:11:GLU:CD	1:A:11:GLU:H	2.12	0.58
1:A:207:VAL:HG21	1:A:270:PHE:CE1	2.39	0.58
1:A:341:LYS:HD3	1:A:341:LYS:O	2.04	0.57
1:A:125:LYS:C	1:A:127:ILE:N	2.61	0.57
1:A:28:TYR:OH	1:A:113:HIS:HD2	1.88	0.57
1:A:341:LYS:O	1:A:343:ASP:N	2.36	0.57
1:A:265:LYS:HB3	1:A:265:LYS:NZ	2.20	0.57
1:A:127:ILE:O	1:A:127:ILE:HD13	2.05	0.56
1:A:6:GLY:HA2	1:A:10:PRO:HG2	1.87	0.56
1:A:182:ALA:HB3	1:A:188:ARG:HG2	1.87	0.56
1:A:246:ILE:HD11	1:A:272:ILE:HD12	1.88	0.56
1:A:299:VAL:HG23	1:A:432:VAL:CG1	2.36	0.55
1:A:378:ILE:HD12	1:A:394:VAL:HB	1.89	0.55
1:A:286:GLU:OE1	1:A:369:ARG:HD3	2.06	0.55
1:A:259:ILE:HG23	1:A:260:PRO:HD2	1.89	0.55
1:A:171:ILE:H	1:A:171:ILE:HD13	1.72	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:158:TYR:HB2	1:A:171:ILE:CD1	2.37	0.55
1:A:364:GLN:O	1:A:365:ASN:OD1	2.24	0.55
1:A:164:ASP:CB	1:A:167:THR:OG1	2.53	0.54
1:A:82:LYS:HD3	1:A:110:ASP:OD1	2.07	0.54
1:A:218:SER:OG	1:A:221:ARG:HB2	2.07	0.54
1:A:125:LYS:O	1:A:127:ILE:N	2.41	0.54
1:A:180:LEU:HB2	1:A:181:ARG:HD2	1.89	0.54
1:A:181:ARG:HD2	1:A:181:ARG:N	2.23	0.54
1:A:127:ILE:HD13	1:A:127:ILE:C	2.33	0.53
1:A:243:ALA:HB2	1:A:269:ARG:NH1	2.22	0.53
1:A:166:ASP:O	1:A:200:ARG:HA	2.08	0.53
1:A:362:ARG:HD3	5:A:844:HOH:O	2.08	0.53
1:A:309:VAL:HG12	1:A:310:GLU:N	2.22	0.53
1:A:158:TYR:O	1:A:171:ILE:HD13	2.08	0.53
1:A:365:ASN:OD1	3:C:1:NAG:H2	2.09	0.53
1:A:340:ASN:CG	1:A:341:LYS:N	2.67	0.53
1:A:288:ARG:HG2	5:A:672:HOH:O	2.09	0.53
1:A:5:GLN:HE21	1:A:5:GLN:HA	1.74	0.52
1:A:144:LYS:HG3	1:A:281:HIS:HD2	1.74	0.52
1:A:365:ASN:OD1	3:C:1:NAG:C2	2.57	0.52
1:A:364:GLN:C	1:A:365:ASN:OD1	2.52	0.52
1:A:6:GLY:HA2	1:A:10:PRO:HG3	1.91	0.51
1:A:140:ASP:OD2	1:A:144:LYS:HB2	2.10	0.51
1:A:366:LEU:HD12	1:A:366:LEU:C	2.35	0.51
1:A:243:ALA:N	1:A:268:CYS:O	2.43	0.51
1:A:264:ALA:HB3	1:A:271:GLU:HB2	1.92	0.51
1:A:304:GLN:N	1:A:305:PRO:HD2	2.26	0.51
1:A:251:PRO:HG2	1:A:255:ARG:HD2	1.92	0.51
1:A:148:LEU:O	1:A:152:ILE:CD1	2.59	0.50
1:A:5:GLN:HA	1:A:5:GLN:NE2	2.26	0.50
1:A:5:GLN:HG3	1:A:6:GLY:N	2.26	0.50
1:A:207:VAL:CG2	1:A:242:PRO:HB3	2.40	0.49
1:A:246:ILE:CD1	1:A:272:ILE:HD12	2.42	0.49
1:A:395:ILE:N	1:A:395:ILE:HD12	2.27	0.49
1:A:76:ASN:HB3	5:A:644:HOH:O	2.12	0.48
1:A:200:ARG:O	1:A:200:ARG:HG2	2.14	0.48
1:A:72:LEU:HD12	1:A:92:ILE:O	2.13	0.48
1:A:80:ASN:ND2	1:A:87:LYS:HZ2	2.12	0.48
1:A:251:PRO:HG3	1:A:275:VAL:HG12	1.93	0.48
1:A:343:ASP:HA	5:A:925:HOH:O	2.13	0.48
1:A:125:LYS:HB2	1:A:127:ILE:HG22	1.96	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:309:VAL:HG13	1:A:408:ARG:HH21	1.76	0.48
1:A:215:LYS:HE2	1:A:223:VAL:CG2	2.44	0.47
1:A:40:GLN:HB3	1:A:57:PHE:HB2	1.96	0.47
1:A:234:PRO:O	1:A:235:ASP:HB2	2.13	0.47
1:A:6:GLY:C	1:A:10:PRO:HG3	2.39	0.47
1:A:152:ILE:HD12	1:A:152:ILE:N	2.30	0.47
1:A:432:VAL:HG12	1:A:432:VAL:OXT	2.14	0.47
1:A:8:GLU:OE2	1:A:68:ALA:CB	2.64	0.46
1:A:259:ILE:CG2	1:A:260:PRO:HD2	2.45	0.46
1:A:5:GLN:HG3	1:A:7:ALA:H	1.80	0.46
1:A:214:LEU:HD21	1:A:222:LEU:HD13	1.98	0.46
1:A:286:GLU:O	1:A:297:HIS:HE1	1.98	0.46
1:A:158:TYR:HB2	1:A:171:ILE:HD11	1.97	0.46
1:A:152:ILE:HG13	1:A:173:VAL:O	2.16	0.46
1:A:237:CYS:O	1:A:238:ASP:HB2	2.14	0.46
1:A:29:LYS:C	1:A:30:ILE:HD12	2.41	0.46
1:A:109:THR:CG2	1:A:110:ASP:N	2.78	0.45
1:A:216:LEU:C	1:A:216:LEU:HD23	2.41	0.45
1:A:30:ILE:HD12	1:A:30:ILE:N	2.30	0.45
1:A:233:GLN:CG	1:A:242:PRO:HD3	2.47	0.45
1:A:243:ALA:HB3	1:A:269:ARG:HD2	1.99	0.45
1:A:255:ARG:NH2	1:A:258:THR:OG1	2.50	0.45
1:A:18:GLU:HA	1:A:28:TYR:O	2.17	0.44
1:A:265:LYS:NZ	1:A:265:LYS:CB	2.80	0.44
1:A:131:ASN:HB2	5:A:762:HOH:O	2.16	0.44
1:A:6:GLY:CA	1:A:10:PRO:HG3	2.48	0.44
1:A:70:ARG:NH1	5:A:855:HOH:O	2.49	0.44
1:A:305:PRO:O	1:A:308:ARG:HG2	2.18	0.44
1:A:293:SER:H	1:A:297:HIS:CD2	2.16	0.44
1:A:181:ARG:N	1:A:181:ARG:CD	2.81	0.44
1:A:109:THR:CG2	1:A:110:ASP:H	2.22	0.43
1:A:158:TYR:HB2	1:A:171:ILE:HD13	2.00	0.43
1:A:149:ASN:C	1:A:152:ILE:HD13	2.43	0.43
1:A:265:LYS:HB3	1:A:265:LYS:HZ3	1.84	0.43
1:A:69:SER:HB2	1:A:72:LEU:HB3	1.99	0.43
1:A:270:PHE:H	1:A:270:PHE:HD1	1.65	0.43
1:A:76:ASN:OD1	2:B:1:NAG:C2	2.67	0.43
1:A:432:VAL:CG1	1:A:432:VAL:OXT	2.67	0.43
1:A:270:PHE:CD1	1:A:270:PHE:N	2.87	0.43
1:A:221:ARG:NH1	1:A:247:THR:HG21	2.33	0.42
1:A:185:PRO:HA	1:A:188:ARG:NE	2.29	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:17:TRP:HH2	1:A:100:LEU:HD13	1.85	0.42
1:A:293:SER:N	1:A:297:HIS:HD2	2.05	0.42
1:A:297:HIS:O	1:A:298:ASP:HB2	2.20	0.42
1:A:310:GLU:O	1:A:312:SER:N	2.47	0.42
1:A:160:VAL:HG22	1:A:171:ILE:HD12	1.97	0.42
1:A:366:LEU:HD13	1:A:377:LEU:HD11	2.01	0.42
1:A:340:ASN:CB	5:A:790:HOH:O	2.68	0.41
1:A:189:VAL:HG12	1:A:189:VAL:O	2.20	0.41
1:A:164:ASP:HA	1:A:165:PRO:HD3	1.88	0.41
1:A:10:PRO:HD2	1:A:11:GLU:OE2	2.21	0.41
1:A:162:ASP:HB3	1:A:262:LEU:HD12	2.02	0.41
1:A:343:ASP:N	5:A:925:HOH:O	2.53	0.41
1:A:38:VAL:HB	1:A:41:CYS:SG	2.60	0.41
1:A:92:ILE:HD13	1:A:115:PHE:CG	2.55	0.41
1:A:344:ALA:HB2	1:A:359:VAL:HG22	2.03	0.41
1:A:102:THR:HG23	5:A:635:HOH:O	2.20	0.41
1:A:182:ALA:O	1:A:188:ARG:HD3	2.21	0.40
1:A:378:ILE:HD13	1:A:394:VAL:HA	2.04	0.40
1:A:328:LEU:HD12	1:A:328:LEU:N	2.37	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	426/432 (99%)	392 (92%)	29 (7%)	5 (1%)	10 3

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	340	ASN
1	A	342	LYS

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Mol	Chain	Res	Type
1	A	201	GLY
1	A	126	ASN
1	A	8	GLU

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	368/370 (100%)	349 (95%)	19 (5%)	21 9

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

Mol	Chain	Res	Type
1	A	11	GLU
1	A	40	GLN
1	A	100	LEU
1	A	127	ILE
1	A	148	LEU
1	A	152	ILE
1	A	164	ASP
1	A	168	SER
1	A	169	LEU
1	A	171	ILE
1	A	212	GLU
1	A	255	ARG
1	A	265	LYS
1	A	313	ASP
1	A	330	ILE
1	A	364	GLN
1	A	366	LEU
1	A	394	VAL
1	A	398	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (13) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	5	GLN
1	A	25	ASN
1	A	35	ASN
1	A	40	GLN
1	A	58	HIS
1	A	80	ASN
1	A	85	ASN
1	A	113	HIS
1	A	233	GLN
1	A	297	HIS
1	A	304	GLN
1	A	364	GLN
1	A	401	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

7 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	2,1	14,14,15	1.36	2 (14%)	17,19,21	1.37	1 (5%)
2	NAG	B	2	2	14,14,15	0.54	0	17,19,21	0.93	0
2	BMA	B	3	2	11,11,12	0.59	0	15,15,17	0.39	0
2	MAN	B	4	2	11,11,12	0.50	0	15,15,17	0.50	0
2	MAN	B	5	2	11,11,12	0.43	0	15,15,17	0.59	0
3	NAG	C	1	1,3	14,14,15	1.03	1 (7%)	17,19,21	1.23	2 (11%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	NAG	C	2	3	14,14,15	0.58	0	17,19,21	0.74	1 (5%)

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	2,1	-	0/6/23/26	0/1/1/1
2	NAG	B	2	2	-	2/6/23/26	0/1/1/1
2	BMA	B	3	2	-	0/2/19/22	0/1/1/1
2	MAN	B	4	2	-	0/2/19/22	0/1/1/1
2	MAN	B	5	2	-	2/2/19/22	0/1/1/1
3	NAG	C	1	1,3	-	3/6/23/26	0/1/1/1
3	NAG	C	2	3	-	2/6/23/26	0/1/1/1

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	1	NAG	C1-C2	3.54	1.57	1.52
2	B	1	NAG	C4-C3	2.38	1.58	1.52
3	C	1	NAG	C4-C5	2.04	1.57	1.53

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	1	NAG	O5-C1-C2	-3.38	106.06	111.29
3	C	1	NAG	O5-C1-C2	-3.01	106.63	111.29
3	C	1	NAG	C3-C4-C5	2.25	114.32	110.23
3	C	2	NAG	C2-N2-C7	-2.04	120.16	122.90

There are no chirality outliers.

All (9) 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
2	B	5	MAN	O5-C5-C6-O6
3	C	1	NAG	C8-C7-N2-C2
2	B	2	NAG	C8-C7-N2-C2

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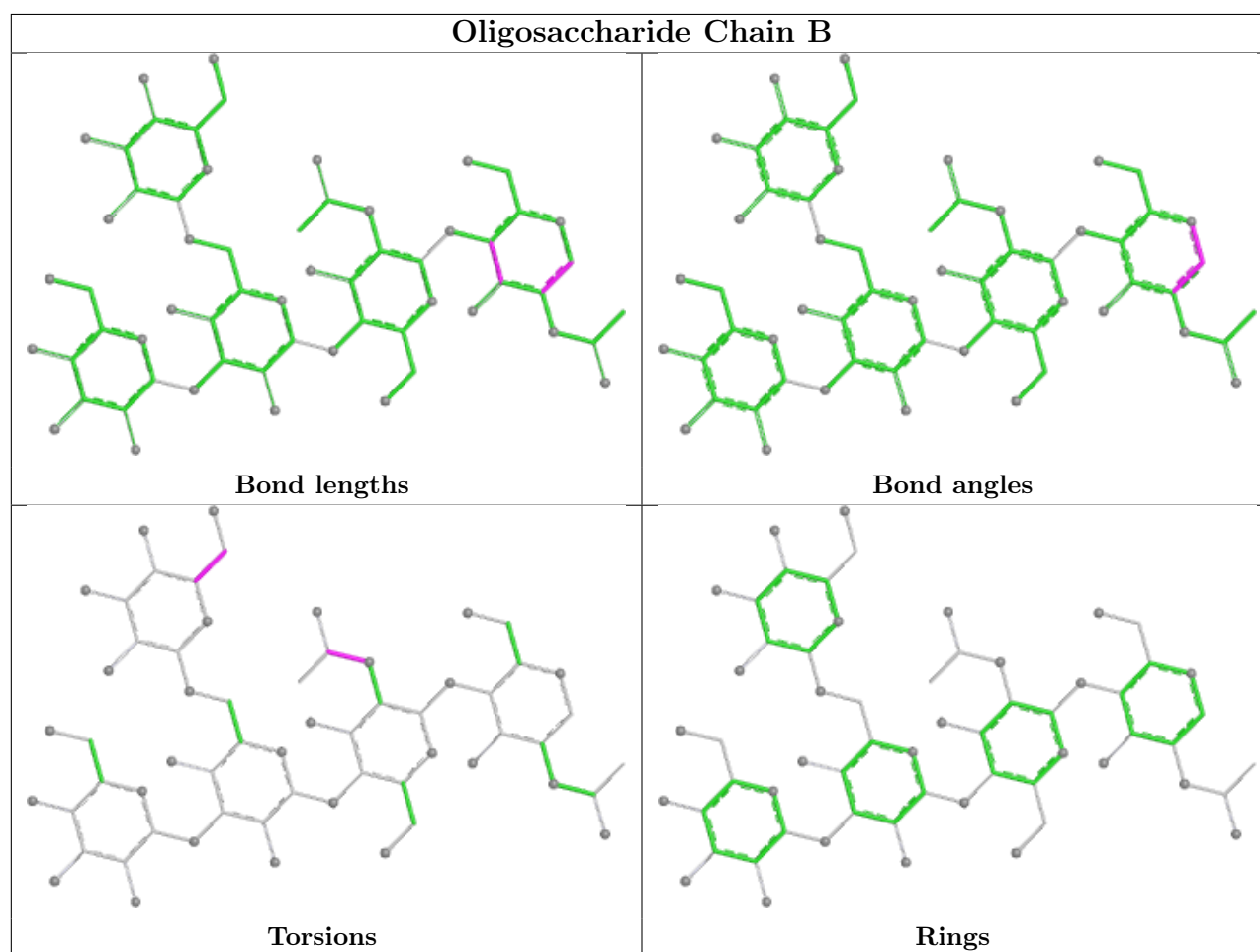
Mol	Chain	Res	Type	Atoms
2	B	2	NAG	O7-C7-N2-C2
3	C	1	NAG	O7-C7-N2-C2
3	C	1	NAG	O5-C5-C6-O6
2	B	5	MAN	C4-C5-C6-O6

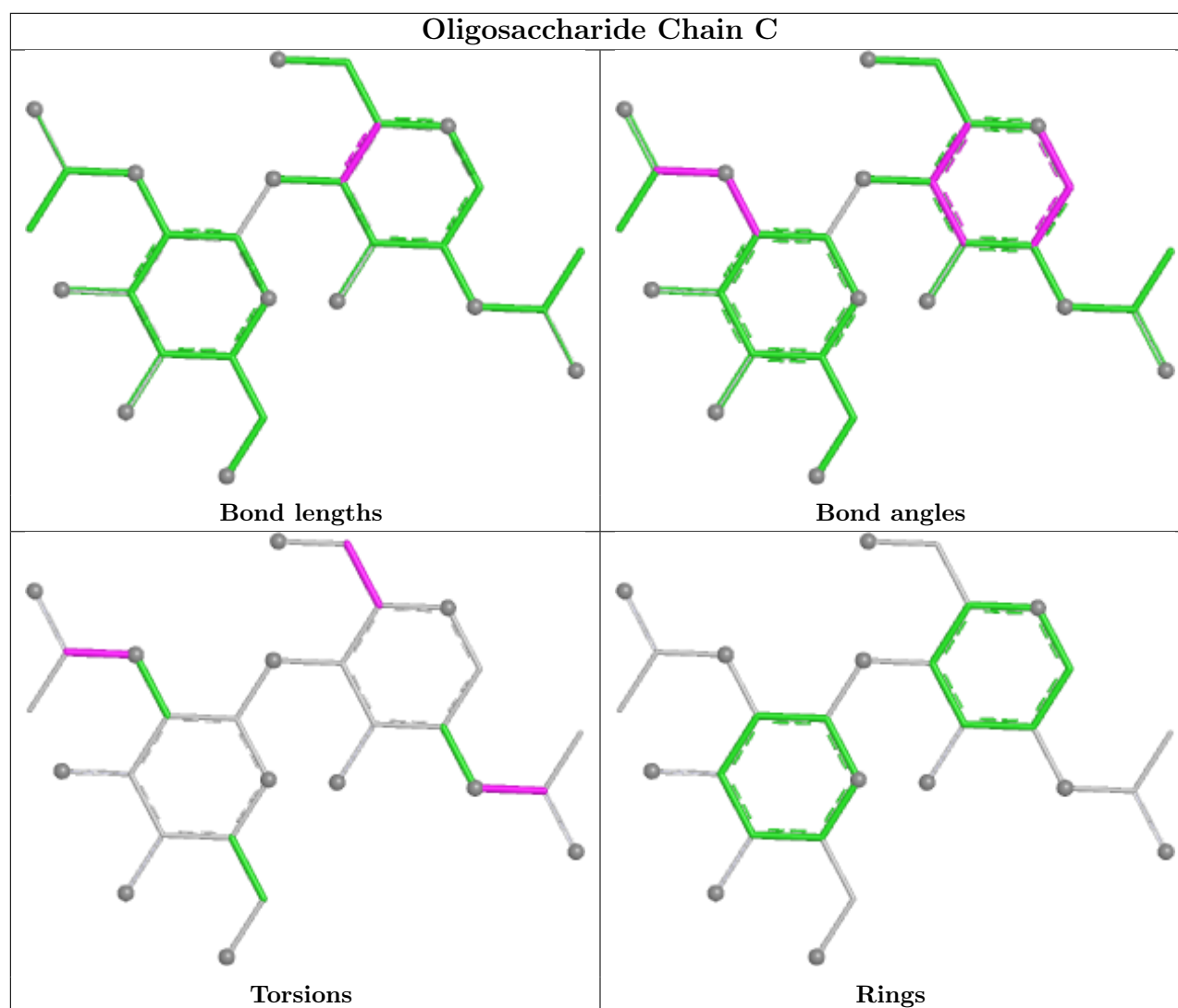
There are no ring outliers.

2 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	1	NAG	1	0
3	C	1	NAG	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 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	$\# Z > 2$
4	GOL	A	4005	-	5,5,5	4.87	5 (100%)	5,5,5	6.09	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
4	GOL	A	4005	-	-	2/4/4/4	-

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	4005	GOL	C3-C2	-8.17	1.20	1.51
4	A	4005	GOL	O1-C1	4.67	1.62	1.42
4	A	4005	GOL	O3-C3	3.44	1.56	1.42
4	A	4005	GOL	C1-C2	-3.12	1.39	1.51
4	A	4005	GOL	O2-C2	-2.89	1.35	1.43

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	4005	GOL	O3-C3-C2	11.06	160.18	110.38
4	A	4005	GOL	O2-C2-C3	7.17	138.86	109.18
4	A	4005	GOL	O1-C1-C2	3.38	125.59	110.38

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	4005	GOL	O1-C1-C2-C3
4	A	4005	GOL	C1-C2-C3-O3

There are no ring outliers.

No monomer is involved in short contacts.

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, 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	428/432 (99%)	0.92	81 (18%) 3 2	13, 28, 68, 83	0

All (81) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	268	CYS	8.3
1	A	237	CYS	8.1
1	A	166	ASP	6.8
1	A	236	PHE	5.7
1	A	182	ALA	5.5
1	A	6	GLY	4.6
1	A	7	ALA	4.6
1	A	165	PRO	4.6
1	A	34	GLY	4.6
1	A	264	ALA	4.5
1	A	340	ASN	4.4
1	A	127	ILE	4.1
1	A	180	LEU	4.1
1	A	309	VAL	4.0
1	A	164	ASP	4.0
1	A	5	GLN	3.9
1	A	341	LYS	3.9
1	A	267	ASN	3.9
1	A	269	ARG	3.8
1	A	183	SER	3.7
1	A	32	ILE	3.6
1	A	315	LEU	3.4
1	A	201	GLY	3.3
1	A	181	ARG	3.3
1	A	41	CYS	3.3
1	A	185	PRO	3.3
1	A	168	SER	3.2

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Mol	Chain	Res	Type	RSRZ
1	A	125	LYS	3.2
1	A	238	ASP	3.2
1	A	33	CYS	3.1
1	A	240	HIS	3.0
1	A	308	ARG	3.0
1	A	239	GLY	3.0
1	A	403	ALA	3.0
1	A	184	SER	2.9
1	A	235	ASP	2.9
1	A	191	PRO	2.9
1	A	266	SER	2.9
1	A	312	SER	2.9
1	A	8	GLU	2.8
1	A	231	ALA	2.8
1	A	163	SER	2.7
1	A	40	GLN	2.7
1	A	233	GLN	2.7
1	A	241	SER	2.7
1	A	179	VAL	2.6
1	A	70	ARG	2.6
1	A	76	ASN	2.6
1	A	218	SER	2.6
1	A	219	ASN	2.6
1	A	311	ALA	2.5
1	A	38	VAL	2.5
1	A	338	ILE	2.4
1	A	203	ARG	2.4
1	A	199	VAL	2.4
1	A	205	PHE	2.4
1	A	343	ASP	2.4
1	A	187	VAL	2.4
1	A	10	PRO	2.3
1	A	234	PRO	2.3
1	A	227	VAL	2.3
1	A	432	VAL	2.3
1	A	39	ALA	2.2
1	A	378	ILE	2.2
1	A	339	CYS	2.2
1	A	192	THR	2.2
1	A	281	HIS	2.2
1	A	211	GLN	2.2
1	A	365	ASN	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	406	ASN	2.1
1	A	171	ILE	2.1
1	A	352	ALA	2.1
1	A	405	ASN	2.1
1	A	212	GLU	2.1
1	A	186	GLN	2.1
1	A	188	ARG	2.1
1	A	265	LYS	2.0
1	A	245	THR	2.0
1	A	202	ASP	2.0
1	A	342	LYS	2.0
1	A	167	THR	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](#)

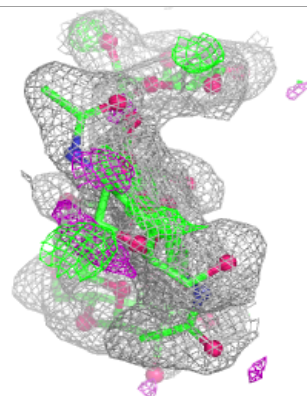
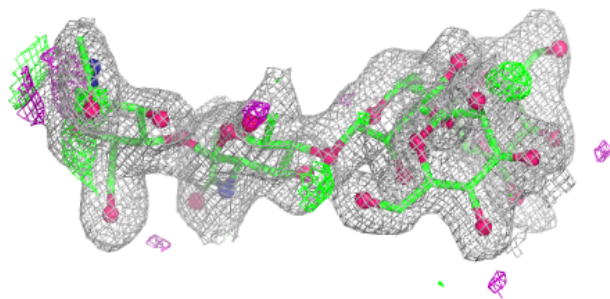
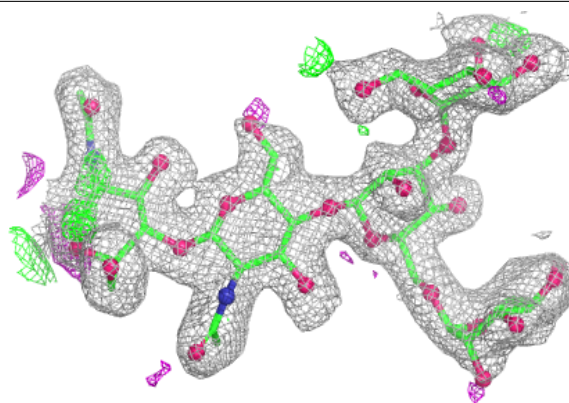
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	NAG	C	2	14/15	0.59	0.18	60,62,63,64	0
3	NAG	C	1	14/15	0.62	0.19	47,51,55,57	0
2	NAG	B	1	14/15	0.80	0.14	27,30,32,33	0
2	NAG	B	2	14/15	0.80	0.11	30,33,37,39	0
2	MAN	B	5	11/12	0.81	0.12	40,42,45,47	0
2	BMA	B	3	11/12	0.91	0.09	22,27,31,36	0
2	MAN	B	4	11/12	0.92	0.09	19,21,25,31	0

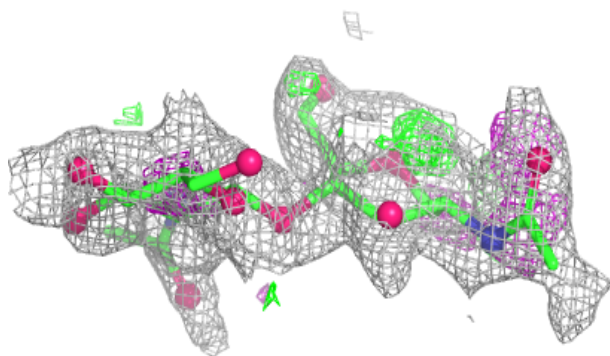
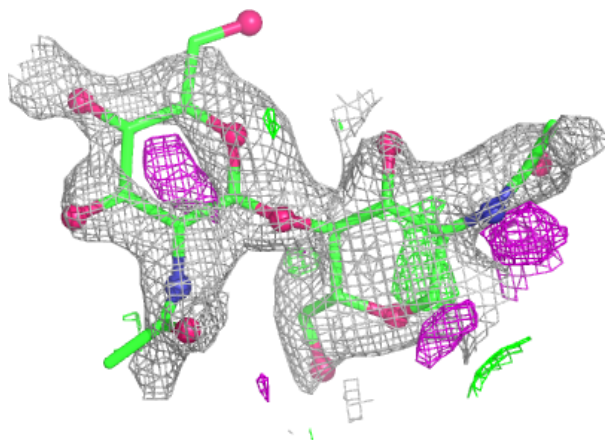
The following is a graphical depiction of the model fit to experimental electron density for oligosaccharide. Each fit is shown from different orientation to approximate a three-dimensional view.

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)

**Electron density around Chain C:**

$2mF_o-DF_c$ (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
4	GOL	A	4005	6/6	0.84	0.12	36,39,41,44	0

6.5 Other polymers [i](#)

There are no such residues in this entry.