



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 12, 2026 – 04:02 AM UTC

PDB ID : 2OLH / pdb_00002olh
Title : Crystal structure of a signalling protein (SPG-40) complex with cellobiose at 2.78 Å resolution
Authors : Sharma, P.; Singh, N.; Sharma, S.; Bhushan, A.; Kaur, P.; Singh, T.P.
Deposited on : 2007-01-19
Resolution : 2.78 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)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

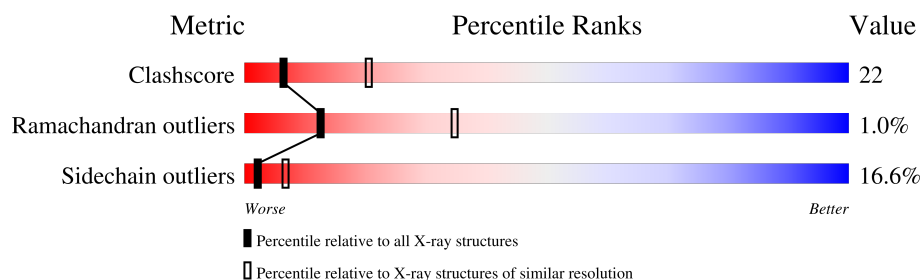
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.78 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))
Clashscore	190562	5693 (2.80-2.76)
Ramachandran outliers	187476	5590 (2.80-2.76)
Sidechain outliers	187428	5592 (2.80-2.76)

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 $\leq 5\%$.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	361	
2	B	3	
3	C	2	

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
2	NAG	B	2	X	-	-	-
2	MAN	B	3	X	-	-	-

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Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	BGC	C	2	X	-	-	-

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 3067 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 Chitinase-3-like protein 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	361	Total	C	N	O	S	0	0	0
			2876	1836	508	523	9			

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	33	VAL	ILE	SEE REMARK 999	UNP Q8SPQ0
A	131	ALA	GLY	SEE REMARK 999	UNP Q8SPQ0
A	205	ASN	GLN	SEE REMARK 999	UNP Q8SPQ0
A	206	SER	GLU	SEE REMARK 999	UNP Q8SPQ0
A	?	-	ASP	SEE REMARK 999	UNP Q8SPQ0
A	361	ARG	GLU	SEE REMARK 999	UNP Q8SPQ0

- Molecule 2 is an oligosaccharide called alpha-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	3	Total	C	N	O	0	0	0
			39	22	2	15			

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



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

- Molecule 4 is water.

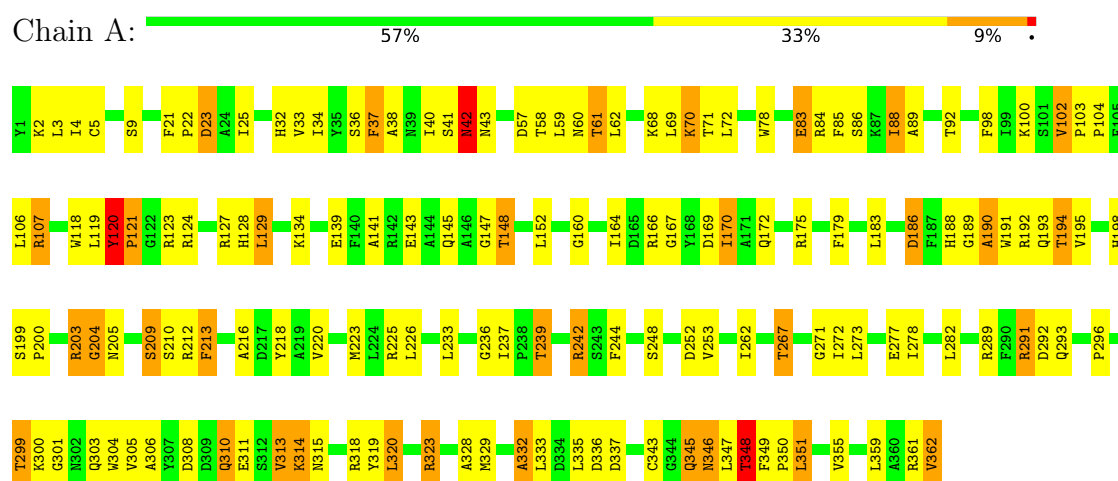
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	129	Total	O	0	0
			129	129		

3 Residue-property plots

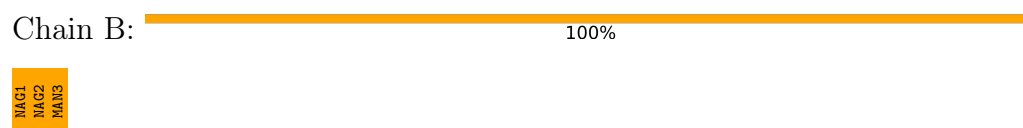
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry. Residues are color-coded 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. 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.

Note EDS was not executed.

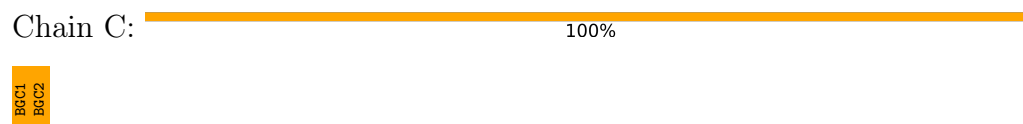
- Molecule 1: Chitinase-3-like protein 1



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



- Molecule 3: beta-D-glucopyranose-(1-4)-beta-D-glucopyranose



4 Data and refinement statistics

Xtriage (Phenix) and EDS were not executed - this section is therefore incomplete.

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	62.69Å 66.45Å 107.53Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	19.94 – 2.78	Depositor
% Data completeness (in resolution range)	97.9 (19.94-2.78)	Depositor
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	CNS 0.9	Depositor
R, R_{free}	0.191 , 0.208	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	3067	wwPDB-VP
Average B, all atoms (Å ²)	33.0	wwPDB-VP

5 Model quality

5.1 Standard geometry

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

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	0.54	0/2952	1.23	32/4001 (0.8%)

There are no bond length outliers.

All (32) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	212	ARG	N-CA-C	11.31	126.22	111.75
1	A	343	CYS	N-CA-C	10.62	124.82	112.93
1	A	36	SER	N-CA-C	10.24	125.25	108.96
1	A	205	ASN	N-CA-C	9.30	123.23	111.24
1	A	188	HIS	N-CA-C	-9.20	96.06	109.59
1	A	209	SER	N-CA-C	8.57	123.83	113.12
1	A	210	SER	CA-C-N	8.38	132.62	120.38
1	A	210	SER	C-N-CA	8.38	132.62	120.38
1	A	190	ALA	N-CA-C	-7.20	102.79	112.94
1	A	37	PHE	N-CA-C	7.01	125.73	110.80
1	A	193	GLN	N-CA-C	-6.90	103.15	112.26
1	A	346	ASN	N-CA-C	6.84	121.38	112.89
1	A	203	ARG	N-CA-C	6.61	118.17	110.97
1	A	61	THR	N-CA-C	-6.27	105.60	113.18
1	A	121	PRO	N-CA-C	-6.27	101.22	111.68
1	A	336	ASP	N-CA-C	-6.10	102.67	110.53
1	A	252	ASP	N-CA-C	5.88	118.47	110.35
1	A	204	GLY	N-CA-C	5.87	122.01	112.61
1	A	191	TRP	N-CA-C	5.78	121.77	114.31
1	A	166	ARG	N-CA-C	5.66	117.12	111.07
1	A	89	ALA	N-CA-C	5.59	117.46	111.36
1	A	40	ILE	N-CA-C	-5.54	100.14	108.12
1	A	42	ASN	N-CA-C	-5.48	105.90	112.58
1	A	62	LEU	N-CA-C	-5.45	105.45	111.71

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	167	GLY	N-CA-C	5.41	126.00	113.18
1	A	120	TYR	CA-C-N	5.36	125.82	120.14
1	A	120	TYR	C-N-CA	5.36	125.82	120.14
1	A	38	ALA	N-CA-C	-5.34	101.57	110.17
1	A	348	THR	N-CA-C	-5.14	101.33	109.76
1	A	199	SER	CA-C-N	-5.12	114.68	119.85
1	A	199	SER	C-N-CA	-5.12	114.68	119.85
1	A	102	VAL	N-CA-C	5.09	119.89	108.88

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	2876	0	2817	124	0
2	B	39	0	34	3	0
3	C	23	0	21	5	0
4	A	129	0	0	27	0
All	All	3067	0	2872	127	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 22.

All (127) 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:78:TRP:HB2	3:C:2:BGC:H6C1	1.44	1.00
1:A:242:ARG:HH22	3:C:1:BGC:H2	1.26	0.97
1:A:78:TRP:HB2	3:C:2:BGC:C6	1.97	0.94
1:A:141:ALA:HB2	4:A:946:HOH:O	1.70	0.90
1:A:203:ARG:HB2	4:A:885:HOH:O	1.79	0.82
2:B:1:NAG:H62	2:B:2:NAG:O7	1.78	0.82
1:A:25:ILE:HD12	1:A:333:LEU:HD13	1.61	0.81
1:A:134:LYS:HE2	4:A:961:HOH:O	1.84	0.77

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:192:ARG:HD3	1:A:194:THR:HG22	1.66	0.76
1:A:57:ASP:O	1:A:61:THR:HG23	1.87	0.73
1:A:242:ARG:NH2	3:C:1:BGC:H2	2.02	0.71
1:A:200:PRO:HB3	1:A:293:GLN:HB3	1.71	0.71
1:A:213:PHE:HB3	1:A:218:TYR:CD2	2.26	0.70
1:A:348:THR:HG23	1:A:349:PHE:HD2	1.57	0.70
1:A:278:ILE:CD1	1:A:305:VAL:HG11	2.21	0.69
1:A:134:LYS:HE3	4:A:866:HOH:O	1.93	0.68
1:A:348:THR:HG23	1:A:349:PHE:CD2	2.29	0.67
1:A:103:PRO:HB2	1:A:104:PRO:HD3	1.76	0.66
1:A:239:THR:HG22	1:A:335:LEU:HB2	1.77	0.66
1:A:278:ILE:HD12	1:A:305:VAL:HG11	1.78	0.65
1:A:267:THR:HB	1:A:277:GLU:OE1	1.96	0.64
1:A:124:ARG:HG2	4:A:941:HOH:O	1.98	0.64
1:A:323:ARG:HH11	1:A:323:ARG:HG2	1.64	0.63
1:A:160:GLY:O	1:A:164:ILE:HG13	1.98	0.62
1:A:84:ARG:HD2	4:A:958:HOH:O	2.00	0.61
4:A:863:HOH:O	2:B:3:MAN:H62	2.00	0.61
1:A:332:ALA:HB3	1:A:335:LEU:HD12	1.83	0.60
1:A:170:ILE:CD1	1:A:223:MET:HE1	2.31	0.60
1:A:169:ASP:OD2	1:A:172:GLN:HG3	2.02	0.60
1:A:291:ARG:HB2	1:A:291:ARG:CZ	2.31	0.60
1:A:345:GLN:HB3	4:A:953:HOH:O	2.04	0.58
1:A:271:GLY:C	1:A:272:ILE:HD12	2.29	0.58
1:A:323:ARG:HG2	1:A:323:ARG:NH1	2.18	0.58
1:A:332:ALA:CB	1:A:335:LEU:HD12	2.33	0.57
1:A:236:GLY:HA3	1:A:329:MET:HE3	1.86	0.57
1:A:124:ARG:HD2	4:A:887:HOH:O	2.05	0.56
1:A:170:ILE:HG13	1:A:223:MET:HE1	1.87	0.56
1:A:107:ARG:HD3	1:A:143:GLU:OE2	2.06	0.56
1:A:107:ARG:NH1	1:A:148:THR:HG21	2.21	0.56
1:A:278:ILE:HD11	1:A:305:VAL:HG11	1.89	0.55
1:A:107:ARG:NH1	1:A:148:THR:CG2	2.70	0.54
1:A:83:GLU:HB3	4:A:855:HOH:O	2.06	0.54
1:A:186:ASP:CG	1:A:186:ASP:O	2.50	0.54
1:A:123:ARG:CD	4:A:922:HOH:O	2.56	0.54
1:A:34:ILE:HG12	1:A:72:LEU:HB2	1.91	0.53
1:A:361:ARG:O	1:A:361:ARG:HG3	2.09	0.53
1:A:127:ARG:HB3	4:A:960:HOH:O	2.07	0.53
1:A:361:ARG:O	1:A:362:VAL:HG12	2.11	0.51
1:A:216:ALA:O	1:A:220:VAL:HG23	2.11	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:332:ALA:HB1	1:A:335:LEU:HG	1.92	0.51
1:A:361:ARG:O	1:A:362:VAL:CB	2.59	0.51
1:A:70:LYS:HE3	4:A:950:HOH:O	2.11	0.50
1:A:37:PHE:HB2	4:A:857:HOH:O	2.12	0.50
1:A:239:THR:O	1:A:239:THR:CG2	2.60	0.50
1:A:78:TRP:CE3	1:A:119:LEU:HD13	2.46	0.50
1:A:204:GLY:HA2	1:A:292:ASP:HB3	1.93	0.50
1:A:318:ARG:HE	1:A:361:ARG:HH12	1.60	0.50
1:A:4:ILE:HD13	1:A:179:PHE:CE2	2.47	0.50
1:A:271:GLY:O	1:A:272:ILE:HD12	2.12	0.50
1:A:5:CYS:HB3	1:A:333:LEU:HG	1.94	0.49
1:A:32:HIS:HA	1:A:70:LYS:O	2.12	0.49
1:A:152:LEU:HD23	4:A:884:HOH:O	2.12	0.49
1:A:273:LEU:HD12	1:A:278:ILE:CD1	2.42	0.49
1:A:78:TRP:HB2	3:C:2:BGC:H6C2	1.91	0.49
1:A:127:ARG:NH1	4:A:878:HOH:O	2.46	0.49
1:A:262:ILE:H	1:A:303:GLN:HE22	1.61	0.49
1:A:2:LYS:HG2	4:A:867:HOH:O	2.13	0.48
1:A:361:ARG:O	1:A:362:VAL:HB	2.13	0.48
1:A:239:THR:HG22	1:A:335:LEU:CB	2.43	0.48
1:A:102:VAL:HB	1:A:103:PRO:HD3	1.95	0.48
1:A:289:ARG:HD3	4:A:964:HOH:O	2.14	0.48
1:A:304:TRP:CZ3	1:A:306:ALA:HB2	2.49	0.48
1:A:319:TYR:CZ	1:A:323:ARG:HD2	2.49	0.47
1:A:192:ARG:HD3	1:A:194:THR:CG2	2.39	0.47
1:A:313:VAL:HG13	1:A:355:VAL:HG22	1.95	0.47
1:A:320:LEU:CD1	1:A:328:ALA:HB2	2.45	0.47
1:A:213:PHE:HB3	1:A:218:TYR:CE2	2.49	0.47
1:A:237:ILE:CG2	1:A:313:VAL:HG22	2.44	0.47
1:A:128:HIS:HE1	4:A:912:HOH:O	1.98	0.46
1:A:253:VAL:HG13	4:A:853:HOH:O	2.14	0.46
1:A:170:ILE:CG1	1:A:223:MET:HE1	2.44	0.46
1:A:323:ARG:HH11	1:A:323:ARG:CG	2.28	0.46
1:A:123:ARG:HD2	4:A:913:HOH:O	2.14	0.46
1:A:320:LEU:HD11	1:A:328:ALA:HB2	1.98	0.46
1:A:41:SER:O	1:A:42:ASN:HB2	2.16	0.45
1:A:123:ARG:HD3	4:A:922:HOH:O	2.15	0.45
1:A:120:TYR:CD2	1:A:120:TYR:N	2.78	0.45
1:A:123:ARG:HD2	4:A:922:HOH:O	2.15	0.45
1:A:43:ASN:OD1	1:A:88:ILE:HD11	2.16	0.45
1:A:273:LEU:HD12	1:A:278:ILE:HD13	1.97	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:310:GLN:O	1:A:314:LYS:CG	2.64	0.45
1:A:60:ASN:ND2	1:A:71:THR:OG1	2.49	0.45
1:A:21:PHE:C	1:A:23:ASP:N	2.73	0.44
1:A:107:ARG:NH1	1:A:143:GLU:OE1	2.51	0.44
1:A:345:GLN:HE21	1:A:345:GLN:HB2	1.61	0.44
1:A:237:ILE:HG21	1:A:313:VAL:HG22	2.00	0.44
1:A:239:THR:O	1:A:239:THR:HG23	2.17	0.43
1:A:313:VAL:HG13	1:A:355:VAL:CG2	2.48	0.43
1:A:169:ASP:CG	1:A:172:GLN:HG3	2.43	0.43
1:A:183:LEU:HD23	1:A:329:MET:HE1	2.00	0.43
1:A:78:TRP:HZ2	4:A:959:HOH:O	2.01	0.43
1:A:22:PRO:HB2	1:A:58:THR:HG22	2.01	0.43
1:A:244:PHE:HD2	1:A:304:TRP:CE3	2.37	0.43
1:A:118:TRP:CE3	1:A:129:LEU:HD23	2.53	0.43
1:A:33:VAL:HG11	1:A:59:LEU:HD11	2.00	0.43
1:A:190:ALA:HA	4:A:858:HOH:O	2.17	0.42
1:A:198:HIS:H	1:A:198:HIS:CD2	2.37	0.42
1:A:299:THR:HA	1:A:303:GLN:O	2.19	0.42
1:A:310:GLN:O	1:A:314:LYS:HG3	2.19	0.42
1:A:119:LEU:HA	1:A:119:LEU:HD23	1.47	0.42
1:A:139:GLU:HA	1:A:139:GLU:OE1	2.20	0.42
1:A:320:LEU:C	1:A:320:LEU:HD23	2.44	0.42
1:A:120:TYR:HA	1:A:121:PRO:HD2	1.73	0.41
1:A:300:LYS:O	1:A:301:GLY:C	2.62	0.41
1:A:351:LEU:HA	1:A:351:LEU:HD12	1.81	0.41
1:A:128:HIS:CE1	4:A:912:HOH:O	2.73	0.41
1:A:43:ASN:O	1:A:98:PHE:HA	2.21	0.41
1:A:43:ASN:OD1	1:A:88:ILE:CD1	2.69	0.41
1:A:291:ARG:H	1:A:291:ARG:HG3	1.67	0.41
1:A:236:GLY:HA3	1:A:329:MET:CE	2.50	0.41
1:A:337:ASP:HB3	1:A:350:PRO:CD	2.51	0.40
1:A:107:ARG:HH11	1:A:143:GLU:CD	2.29	0.40
1:A:145:GLN:C	1:A:147:GLY:H	2.30	0.40
1:A:103:PRO:HB3	1:A:143:GLU:HG3	2.04	0.40
1:A:107:ARG:NH1	1:A:148:THR:HG23	2.37	0.40
1:A:291:ARG:HB3	4:A:918:HOH:O	2.21	0.40
2:B:1:NAG:C6	2:B:2:NAG:O7	2.61	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	286/361 (79%)	257 (90%)	26 (9%)	3 (1%)	12	35

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	189	GLY
1	A	332	ALA
1	A	120	TYR

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	302/302 (100%)	252 (83%)	50 (17%)	2	7

All (50) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	3	LEU
1	A	9	SER
1	A	23	ASP
1	A	42	ASN
1	A	68	LYS
1	A	69	LEU
1	A	70	LYS
1	A	83	GLU

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Mol	Chain	Res	Type
1	A	85	PHE
1	A	86	SER
1	A	88	ILE
1	A	92	THR
1	A	100	LYS
1	A	106	LEU
1	A	107	ARG
1	A	129	LEU
1	A	148	THR
1	A	170	ILE
1	A	175	ARG
1	A	186	ASP
1	A	194	THR
1	A	195	VAL
1	A	209	SER
1	A	213	PHE
1	A	225	ARG
1	A	226	LEU
1	A	233	LEU
1	A	239	THR
1	A	242	ARG
1	A	248	SER
1	A	267	THR
1	A	282	LEU
1	A	291	ARG
1	A	296	PRO
1	A	299	THR
1	A	308	ASP
1	A	310	GLN
1	A	311	GLU
1	A	313	VAL
1	A	314	LYS
1	A	315	ASN
1	A	320	LEU
1	A	323	ARG
1	A	345	GLN
1	A	346	ASN
1	A	347	LEU
1	A	348	THR
1	A	351	LEU
1	A	359	LEU
1	A	362	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	60	ASN
1	A	79	ASN
1	A	303	GLN
1	A	310	GLN
1	A	315	ASN

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 ⓘ

5 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	0.84	0	17,19,21	1.33	1 (5%)
2	NAG	B	2	2	14,14,15	1.46	3 (21%)	17,19,21	2.05	3 (17%)
2	MAN	B	3	2	11,11,12	0.99	0	15,15,17	1.38	2 (13%)
3	BGC	C	1	3	12,12,12	2.42	7 (58%)	17,17,17	2.11	3 (17%)
3	BGC	C	2	3	11,11,12	3.33	5 (45%)	15,15,17	1.91	4 (26%)

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	-	2/6/23/26	0/1/1/1
2	NAG	B	2	2	1/1/5/7	1/6/23/26	0/1/1/1
2	MAN	B	3	2	1/1/4/5	2/2/19/22	0/1/1/1
3	BGC	C	1	3	-	2/2/22/22	0/1/1/1
3	BGC	C	2	3	1/1/4/5	2/2/19/22	0/1/1/1

All (15) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	C	2	BGC	O5-C1	7.51	1.56	1.43
3	C	2	BGC	O5-C5	6.55	1.56	1.43
3	C	1	BGC	C4-C5	4.74	1.63	1.53
2	B	2	NAG	C1-C2	3.74	1.57	1.52
3	C	1	BGC	C1-C2	2.89	1.59	1.52
3	C	2	BGC	O2-C2	2.88	1.49	1.43
3	C	1	BGC	O4-C4	2.78	1.49	1.43
3	C	1	BGC	O5-C1	2.55	1.49	1.42
3	C	1	BGC	C4-C3	2.32	1.58	1.52
2	B	2	NAG	C8-C7	2.27	1.55	1.50
3	C	1	BGC	C3-C2	2.23	1.58	1.52
3	C	2	BGC	C6-C5	2.21	1.59	1.51
3	C	2	BGC	C2-C3	2.17	1.55	1.52
3	C	1	BGC	C6-C5	2.04	1.58	1.51
2	B	2	NAG	C3-C2	2.01	1.56	1.52

All (13) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	1	BGC	O4-C4-C3	-6.90	94.12	110.38
2	B	2	NAG	C1-C2-N2	6.66	120.94	110.43
3	C	2	BGC	C1-O5-C5	4.56	118.29	112.19
2	B	3	MAN	C1-C2-C3	4.26	115.84	109.64
2	B	1	NAG	C4-C3-C2	3.89	116.72	111.02
3	C	2	BGC	C1-C2-C3	3.20	114.31	109.64
3	C	2	BGC	O5-C1-C2	-2.91	103.84	110.79
2	B	2	NAG	C2-N2-C7	2.90	126.78	122.90
2	B	2	NAG	C4-C3-C2	2.88	115.24	111.02
2	B	3	MAN	C1-O5-C5	2.72	115.84	112.19
3	C	2	BGC	C6-C5-C4	2.26	118.58	113.02
3	C	1	BGC	C3-C4-C5	-2.05	106.52	110.23
3	C	1	BGC	O2-C2-C3	2.02	115.13	110.38

All (3) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
2	B	2	NAG	C1
2	B	3	MAN	C1
3	C	2	BGC	C1

All (9) torsion outliers are listed below:

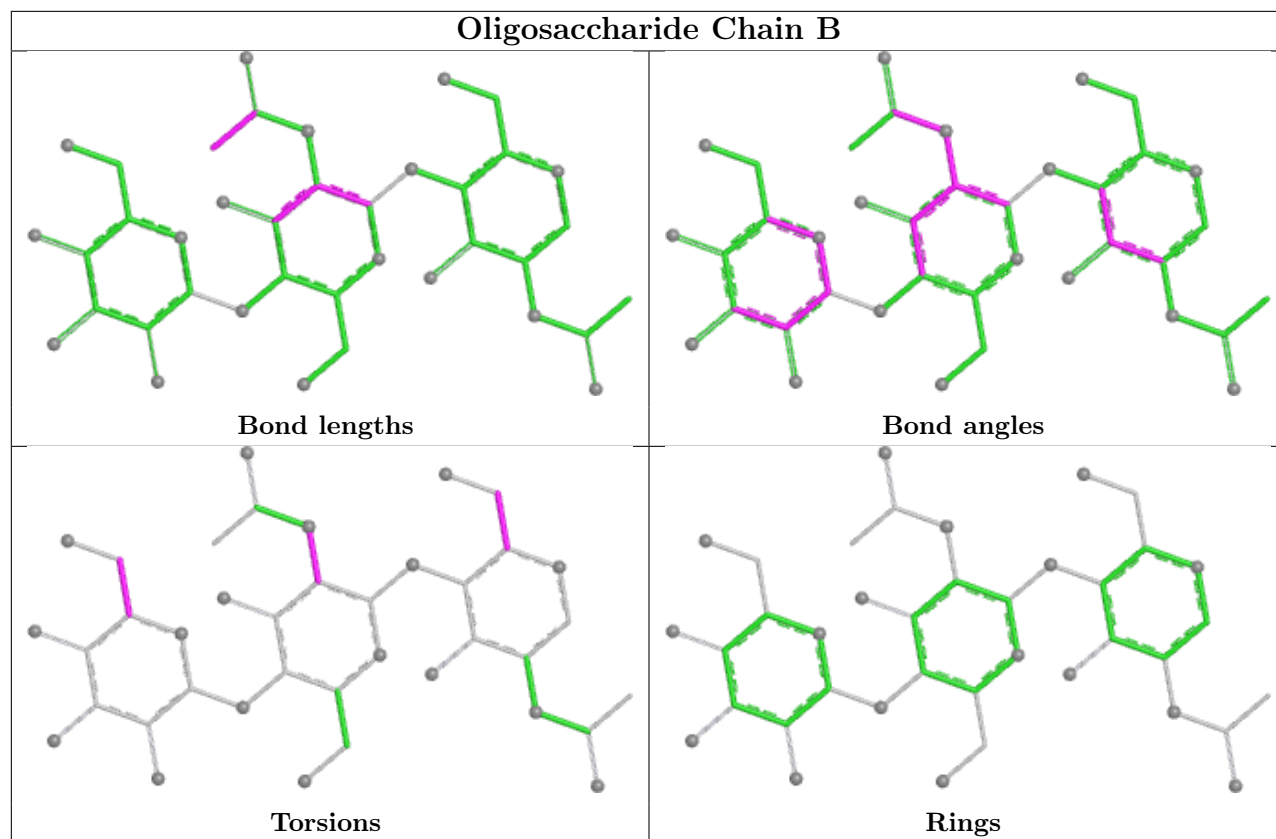
Mol	Chain	Res	Type	Atoms
2	B	2	NAG	C1-C2-N2-C7
3	C	2	BGC	C4-C5-C6-O6
3	C	2	BGC	O5-C5-C6-O6
2	B	3	MAN	C4-C5-C6-O6
3	C	1	BGC	O5-C5-C6-O6
3	C	1	BGC	C4-C5-C6-O6
2	B	3	MAN	O5-C5-C6-O6
2	B	1	NAG	C4-C5-C6-O6
2	B	1	NAG	O5-C5-C6-O6

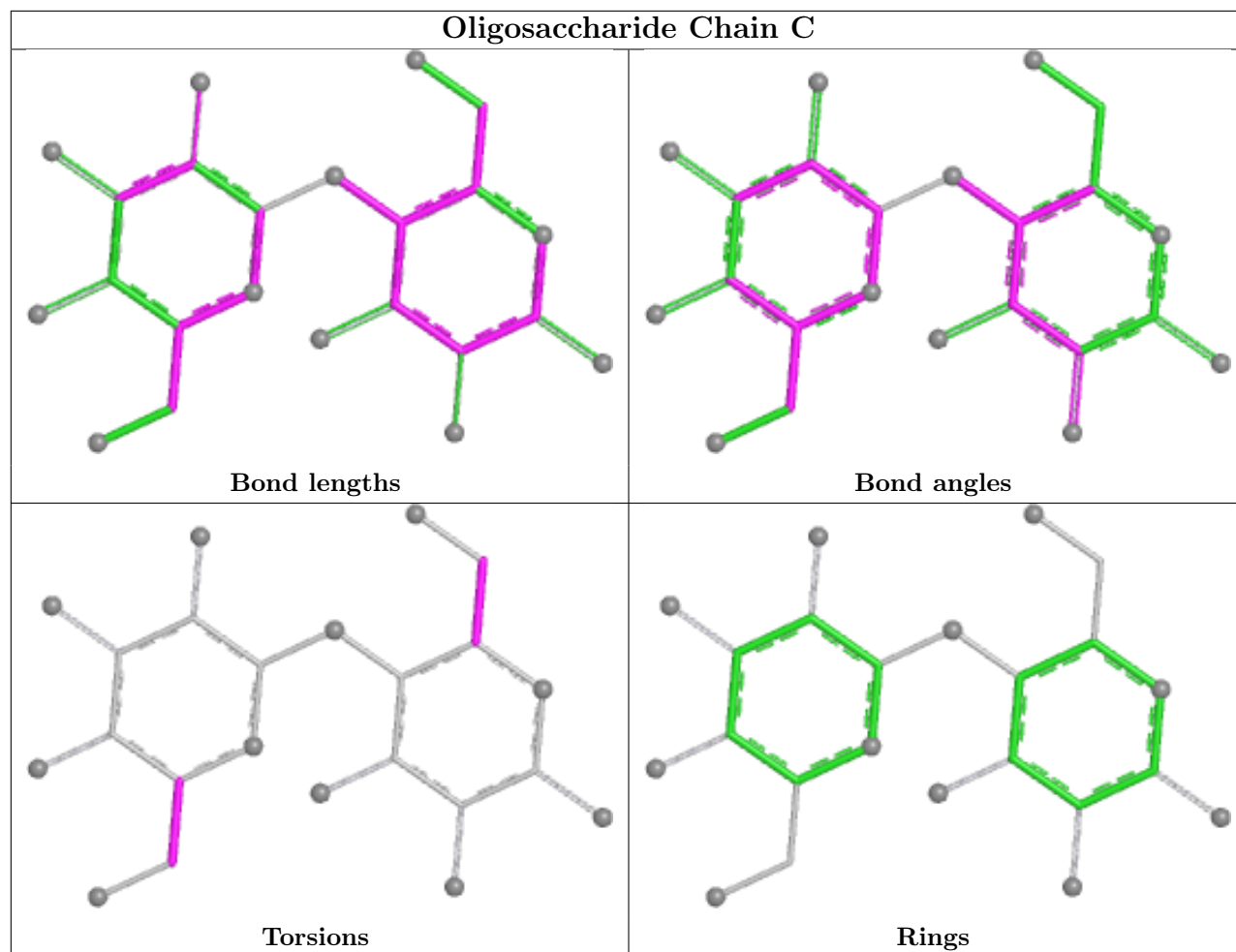
There are no ring outliers.

5 monomers are involved in 8 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	3	MAN	1	0
2	B	1	NAG	2	0
3	C	1	BGC	2	0
2	B	2	NAG	2	0
3	C	2	BGC	3	0

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





5.6 Ligand geometry [i](#)

There are no ligands in this entry.

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

EDS was not executed - this section is therefore empty.

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

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates

EDS was not executed - this section is therefore empty.

6.4 Ligands

EDS was not executed - this section is therefore empty.

6.5 Other polymers

EDS was not executed - this section is therefore empty.