



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 06:19 PM UTC

PDB ID : 6JHX / pdb_00006jhx
Title : Crystal structure of alginate-binding protein AlgQ2 without calcium ion
Authors : Okumura, K.; Maruyama, Y.; Murata, K.; Hashimoto, W.
Deposited on : 2019-02-19
Resolution : 2.20 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

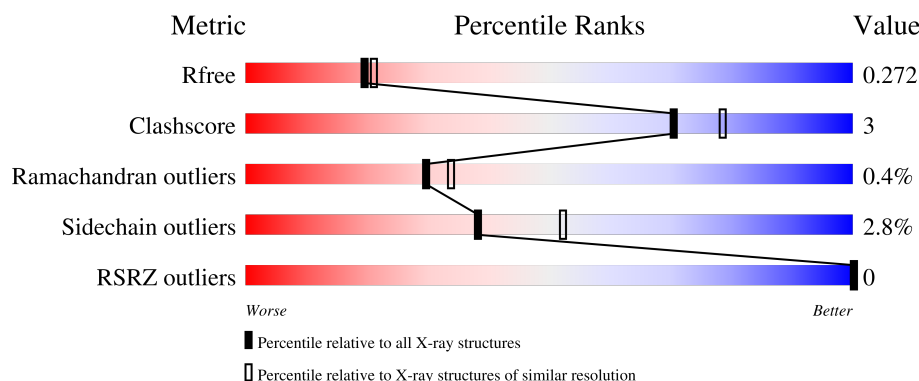
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

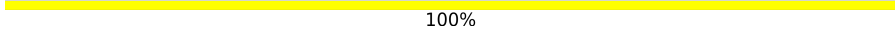
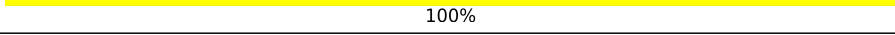
The reported resolution of this entry is 2.20 Å.

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	6164 (2.20-2.20)
Clashscore	190562	6851 (2.20-2.20)
Ramachandran outliers	187476	6768 (2.20-2.20)
Sidechain outliers	187428	6769 (2.20-2.20)
RSRZ outliers	180081	6166 (2.20-2.20)

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	492	 87% 11% .
1	B	492	 89% 9% ..
2	C	3	 100%
2	D	3	 100%

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 8246 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.

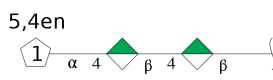
- Molecule 1 is a protein called AlgQ2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	486	Total	C	N	O	S	0	0	0
			4002	2580	684	722	16			
1	B	483	Total	C	N	O	S	0	0	0
			3972	2560	679	717	16			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	179	ALA	ASP	engineered mutation	UNP Q9KWT5
A	180	ALA	GLU	engineered mutation	UNP Q9KWT5
B	179	ALA	ASP	engineered mutation	UNP Q9KWT5
B	180	ALA	GLU	engineered mutation	UNP Q9KWT5

- Molecule 2 is an oligosaccharide called 4-deoxy-alpha-L-erythro-hex-4-enopyranuronic acid-(1-4)-beta-D-mannopyranuronic acid-(1-4)-beta-D-mannopyranuronic acid.



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace
2	C	3	Total	C	O	0	0	0
			36	18	18			
2	D	3	Total	C	O	0	0	0
			36	18	18			

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	109	Total	O	0	0
			109	109		

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	91	Total	O	0	0
			91	91		

4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	45.87Å 60.82Å 87.23Å 80.37° 89.81° 88.17°	Depositor
Resolution (Å)	50.00 – 2.20 50.00 – 2.20	Depositor EDS
% Data completeness (in resolution range)	97.8 (50.00-2.20) 97.8 (50.00-2.20)	Depositor EDS
R_{merge}	0.14	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.92 (at 2.20Å)	Xtriage
Refinement program	REFMAC 5.8.0073	Depositor
R, R_{free}	0.210 , 0.271 0.212 , 0.272	Depositor DCC
R_{free} test set	2350 reflections (4.98%)	wwPDB-VP
Wilson B-factor (Å ²)	27.4	Xtriage
Anisotropy	0.072	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 23.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	0.077 for h,-k,-l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	8246	wwPDB-VP
Average B, all atoms (Å ²)	29.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.56% 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: MAW, BEM

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	0.81	0/4121	0.98	5/5578 (0.1%)
1	B	0.75	0/4090	0.97	2/5537 (0.0%)
All	All	0.78	0/8211	0.98	7/11115 (0.1%)

There are no bond length outliers.

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	437	TYR	N-CA-C	5.84	120.53	113.41
1	A	59	ASN	N-CA-C	5.40	117.59	111.11
1	A	252	ALA	N-CA-C	5.38	117.83	111.33
1	A	205	SER	N-CA-C	5.30	116.06	108.74
1	A	478	VAL	N-CA-C	5.16	115.77	110.36
1	B	252	ALA	CA-C-N	-5.12	114.49	122.73
1	B	252	ALA	C-N-CA	-5.12	114.49	122.73

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	4002	0	3883	33	0
1	B	3972	0	3853	25	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	C	36	0	19	0	0
2	D	36	0	19	0	0
3	A	109	0	0	1	0
3	B	91	0	0	0	0
All	All	8246	0	7774	53	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 3.

All (53) 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:94:LYS:NZ	1:B:477:GLN:HE21	1.64	0.95
1:A:94:LYS:HZ1	1:B:477:GLN:HE21	1.14	0.94
1:A:231:MET:HE1	1:A:453:VAL:HG23	1.80	0.64
1:A:481:MET:HA	1:A:481:MET:HE2	1.81	0.63
1:A:94:LYS:NZ	1:B:477:GLN:NE2	2.45	0.58
1:B:157:ILE:HD13	1:B:197:PHE:HB3	1.85	0.58
1:A:157:ILE:HD13	1:A:197:PHE:HB3	1.87	0.57
1:B:481:MET:HA	1:B:481:MET:HE2	1.86	0.57
1:B:378:LEU:HB3	1:B:383:ALA:HB3	1.89	0.54
1:A:353:THR:HG21	1:A:381:MET:SD	2.48	0.54
1:A:16:HIS:CG	1:A:61:MET:HE2	2.44	0.53
1:B:269:ASP:HB3	1:B:274:THR:OG1	2.09	0.52
1:B:121:ASP:OD1	1:B:123:ASN:HB2	2.10	0.52
1:A:171:ASP:CG	1:A:171:ASP:O	2.54	0.51
1:A:437:TYR:O	1:A:441:VAL:HG23	2.11	0.50
1:A:16:HIS:ND1	1:A:61:MET:HE2	2.27	0.49
1:A:42:LYS:HE3	3:A:658:HOH:O	2.11	0.49
1:B:206:ASP:HB3	1:B:310:GLN:HG3	1.94	0.48
1:A:94:LYS:HZ1	1:B:477:GLN:NE2	1.96	0.48
1:B:188:PRO:O	1:B:191:VAL:HG12	2.14	0.47
1:A:269:ASP:HB3	1:A:274:THR:OG1	2.14	0.47
1:B:231:MET:HE1	1:B:453:VAL:HG23	1.97	0.46
1:A:188:PRO:HB3	1:A:248:PHE:HA	1.99	0.45
1:A:17:MET:O	1:A:24:VAL:HA	2.17	0.45
1:A:201:ARG:NH2	1:A:204:GLY:HA2	2.31	0.45
1:A:121:ASP:OD1	1:A:123:ASN:HB2	2.16	0.45
1:B:360:LYS:HB2	1:B:360:LYS:HE3	1.74	0.45
1:A:106:PHE:CD1	1:A:106:PHE:C	2.96	0.43
1:A:173:ASN:HD21	1:A:179:ALA:HB3	1.84	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:78:ASP:OD1	1:B:431:ARG:NH2	2.51	0.43
1:B:430:GLU:OE1	1:B:484:GLN:HB3	2.17	0.43
1:A:346:ASN:HA	1:A:384:GLN:NE2	2.34	0.43
1:A:113:VAL:HG13	1:A:312:VAL:HB	2.01	0.43
1:A:58:PHE:HA	1:A:61:MET:HE3	2.00	0.43
1:A:62:MET:HE2	1:A:83:TYR:CD2	2.54	0.42
1:B:171:ASP:O	1:B:173:ASN:N	2.52	0.42
1:A:94:LYS:HZ3	1:B:477:GLN:HE21	1.55	0.42
1:A:206:ASP:HB2	1:A:308:SER:O	2.19	0.42
1:A:58:PHE:CD1	1:A:58:PHE:C	2.97	0.42
1:B:188:PRO:HB3	1:B:248:PHE:HA	2.02	0.42
1:B:308:SER:HB3	1:B:417:VAL:CG1	2.49	0.41
1:B:429:GLU:O	1:B:433:ILE:HD13	2.20	0.41
1:B:206:ASP:HB2	1:B:308:SER:O	2.21	0.41
1:B:113:VAL:HG13	1:B:312:VAL:HB	2.02	0.41
1:A:76:LEU:HA	1:A:76:LEU:HD23	1.79	0.41
1:B:309:ARG:HE	1:B:309:ARG:HB2	1.76	0.41
1:A:206:ASP:HB3	1:A:310:GLN:HG3	2.02	0.41
1:A:243:ILE:O	1:A:244:ASP:C	2.64	0.41
1:B:72:GLY:HA2	1:B:316:GLY:O	2.21	0.41
1:B:96:ILE:HA	1:B:100:ALA:HB3	2.02	0.41
1:A:17:MET:HG3	1:A:72:GLY:O	2.21	0.41
1:A:72:GLY:HA2	1:A:316:GLY:O	2.20	0.40
1:A:231:MET:HE1	1:A:453:VAL:CG2	2.50	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	482/492 (98%)	470 (98%)	11 (2%)	1 (0%)	43 51

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	B	479/492 (97%)	465 (97%)	11 (2%)	3 (1%)	21	23
All	All	961/984 (98%)	935 (97%)	22 (2%)	4 (0%)	30	34

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	251	LYS
1	B	172	PRO
1	A	182	PRO
1	B	182	PRO

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	419/422 (99%)	408 (97%)	11 (3%)	40	55
1	B	416/422 (99%)	404 (97%)	12 (3%)	37	51
All	All	835/844 (99%)	812 (97%)	23 (3%)	38	52

All (23) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	49	LYS
1	A	60	LEU
1	A	61	MET
1	A	129	TYR
1	A	135	VAL
1	A	300	LYS
1	A	309	ARG
1	A	395	TYR
1	A	413	LYS
1	A	472	LEU
1	A	477	GLN
1	B	60	LEU

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Mol	Chain	Res	Type
1	B	109	SER
1	B	129	TYR
1	B	184	ILE
1	B	309	ARG
1	B	360	LYS
1	B	395	TYR
1	B	413	LYS
1	B	424	VAL
1	B	425	ASN
1	B	477	GLN
1	B	484	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (12) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	85	GLN
1	A	98	GLN
1	A	233	HIS
1	A	236	GLN
1	A	376	ASN
1	A	425	ASN
1	B	85	GLN
1	B	173	ASN
1	B	257	GLN
1	B	376	ASN
1	B	477	GLN
1	B	480	GLN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

6 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	BEM	C	1	2	13,13,13	1.04	0	18,19,19	2.01	3 (16%)
2	BEM	C	2	2	12,12,13	1.33	2 (16%)	14,17,19	2.06	4 (28%)
2	MAW	C	3	2	10,11,12	1.84	2 (20%)	12,15,17	1.18	2 (16%)
2	BEM	D	1	2	13,13,13	1.01	0	18,19,19	2.14	6 (33%)
2	BEM	D	2	2	12,12,13	1.26	2 (16%)	14,17,19	2.82	5 (35%)
2	MAW	D	3	2	10,11,12	1.94	3 (30%)	12,15,17	1.93	4 (33%)

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	BEM	C	1	2	-	0/4/24/24	0/1/1/1
2	BEM	C	2	2	-	0/4/21/24	0/1/1/1
2	MAW	C	3	2	-	0/4/17/20	0/1/1/1
2	BEM	D	1	2	-	3/4/24/24	0/1/1/1
2	BEM	D	2	2	-	0/4/21/24	0/1/1/1
2	MAW	D	3	2	-	0/4/17/20	0/1/1/1

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	3	MAW	O5-C5	4.52	1.43	1.37
2	D	3	MAW	O5-C5	4.23	1.43	1.37
2	C	3	MAW	C3-C4	2.56	1.53	1.50
2	D	3	MAW	O6B-C6	-2.55	1.23	1.30
2	C	2	BEM	C2-C3	-2.39	1.48	1.52
2	D	2	BEM	O6B-C6	-2.25	1.23	1.30
2	C	2	BEM	O6B-C6	-2.24	1.23	1.30
2	D	3	MAW	C3-C4	2.24	1.53	1.50
2	D	2	BEM	C5-C6	-2.14	1.48	1.53

All (24) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	2	BEM	C1-C2-C3	-7.56	98.63	109.64
2	C	1	BEM	O4-C4-C5	-5.67	96.81	109.76
2	D	1	BEM	C1-O5-C5	-5.55	104.05	112.22
2	D	2	BEM	O2-C2-C3	-4.77	100.28	110.15
2	C	1	BEM	O5-C5-C6	4.20	119.27	106.14
2	C	2	BEM	O2-C2-C3	-4.14	101.58	110.15
2	D	3	MAW	O5-C5-C4	-3.99	121.27	124.94
2	D	3	MAW	O3-C3-C2	-3.66	102.92	109.44
2	D	1	BEM	O5-C5-C6	3.44	116.91	106.14
2	D	1	BEM	O4-C4-C3	3.17	117.85	110.38
2	D	2	BEM	O4-C4-C5	-3.00	102.92	109.76
2	C	2	BEM	O3-C3-C2	-2.99	103.95	110.05
2	D	1	BEM	C3-C4-C5	2.94	114.36	109.30
2	C	3	MAW	O5-C5-C4	-2.93	122.25	124.94
2	C	2	BEM	C2-C3-C4	-2.87	105.81	110.86
2	C	2	BEM	O3-C3-C4	-2.70	104.01	110.38
2	D	2	BEM	O3-C3-C2	-2.41	105.13	110.05
2	D	3	MAW	O6B-C6-C5	2.39	119.48	114.06
2	D	2	BEM	O2-C2-C1	2.37	114.65	109.22
2	C	3	MAW	C1-C2-C3	-2.30	106.29	109.64
2	D	1	BEM	C1-C2-C3	2.30	115.04	110.36
2	C	1	BEM	O4-C4-C3	-2.21	105.18	110.38
2	D	3	MAW	C1-C2-C3	-2.08	106.61	109.64
2	D	1	BEM	O2-C2-C3	-2.00	105.66	110.38

There are no chirality outliers.

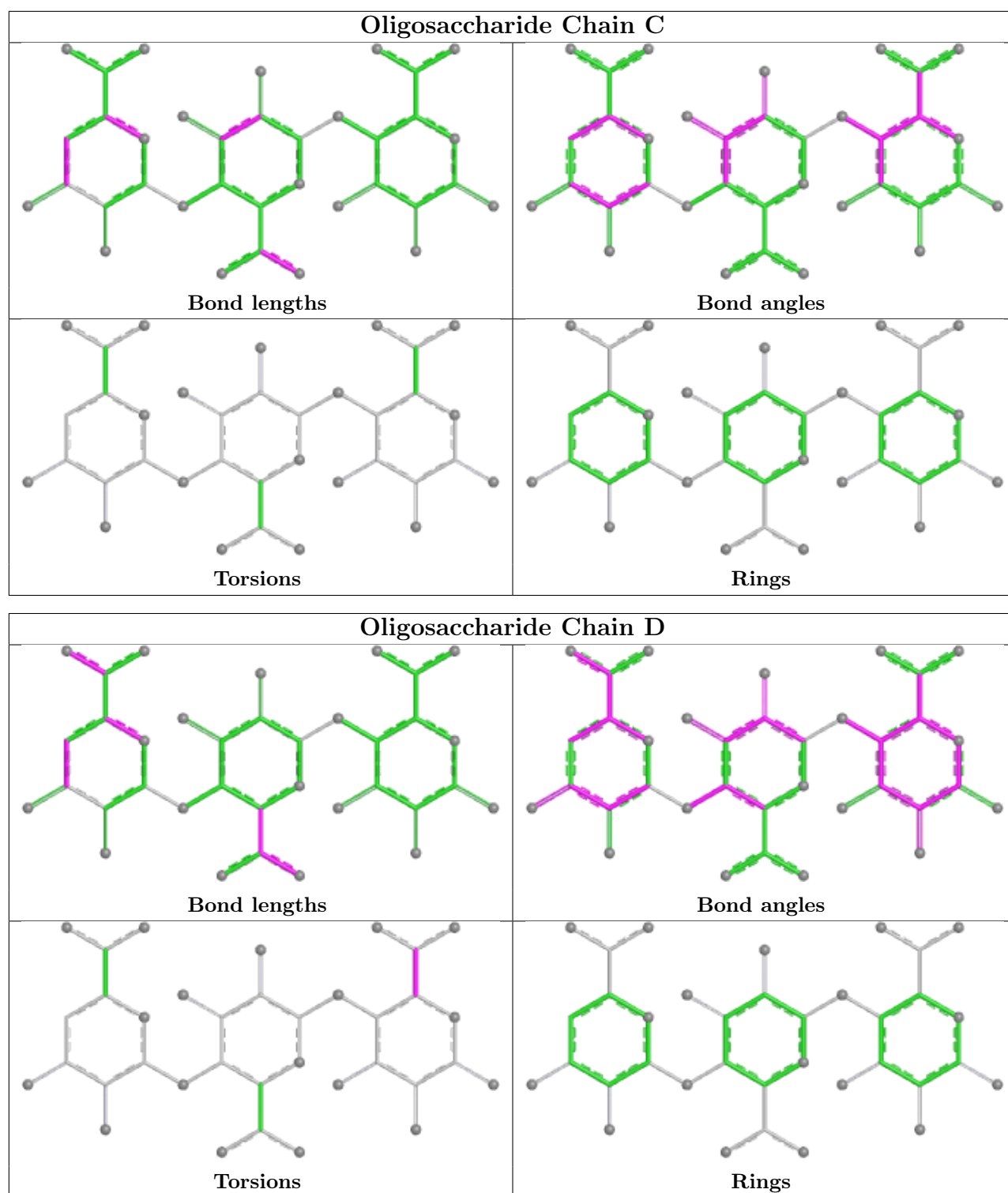
All (3) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	D	1	BEM	C4-C5-C6-O6B
2	D	1	BEM	C4-C5-C6-O6A
2	D	1	BEM	O5-C5-C6-O6B

There are no ring outliers.

No monomer is involved in short contacts.

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



5.6 Ligand geometry ⓘ

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 [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	486/492 (98%)	-1.34	0 100 100	16, 27, 43, 68	0
1	B	483/492 (98%)	-1.34	0 100 100	18, 29, 45, 72	0
All	All	969/984 (98%)	-1.34	0 100 100	16, 28, 44, 72	0

There are no RSRZ outliers to report.

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](#)

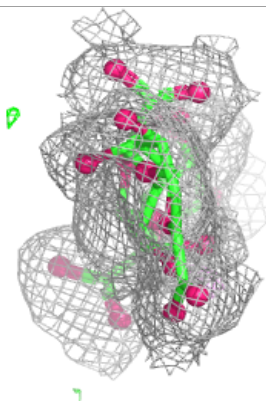
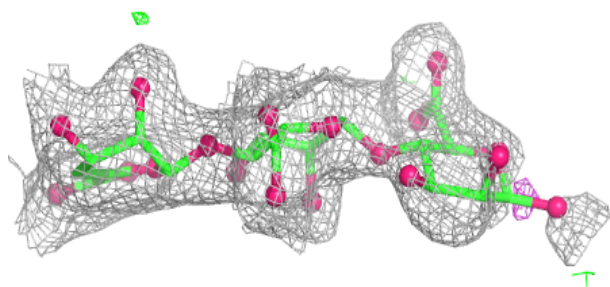
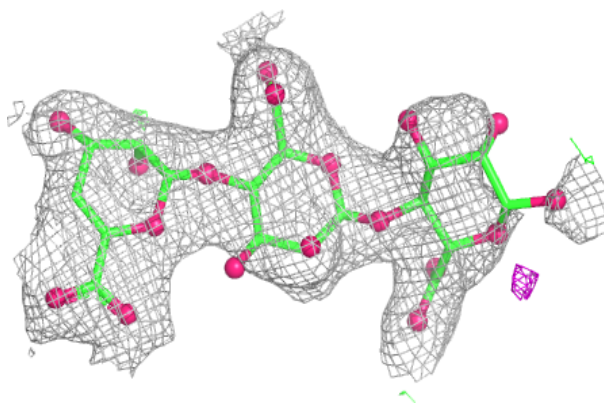
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled ‘Q< 0.9’ lists 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	BEM	D	1	13/13	0.98	0.04	30,33,37,40	0
2	BEM	C	2	12/13	0.99	0.03	27,33,36,46	0
2	MAW	C	3	11/12	0.99	0.02	22,26,28,29	0
2	BEM	C	1	13/13	0.99	0.04	33,42,55,55	0
2	BEM	D	2	12/13	0.99	0.03	23,31,37,37	0
2	MAW	D	3	11/12	0.99	0.02	21,22,23,25	0

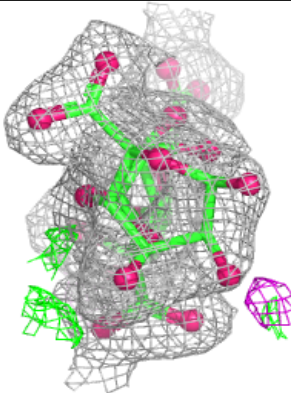
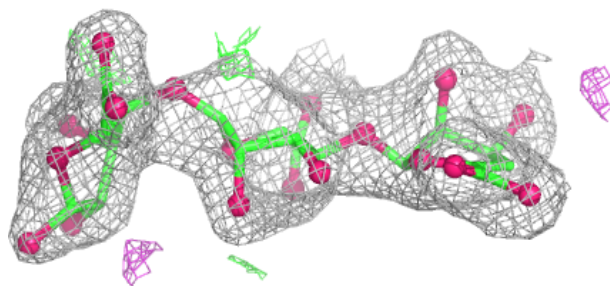
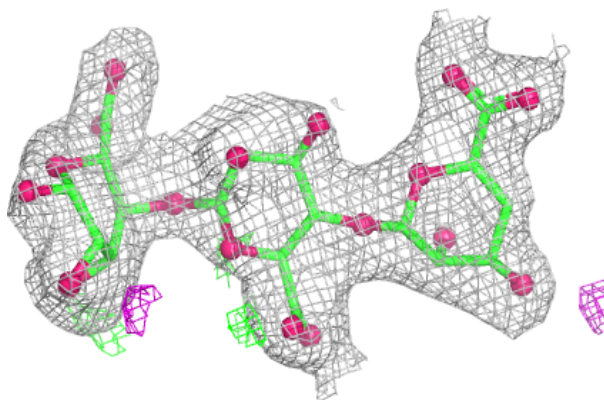
The following is a graphical depiction of the model fit to experimental electron density for oligosaccharide. Each fit is shown from different orientation to approximate a three-dimensional view.

Electron density around Chain C:

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

**Electron density around Chain D:**

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



6.4 Ligands

There are no ligands in this entry.

6.5 Other polymers

There are no such residues in this entry.