



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 05:55 AM UTC

PDB ID : 4X9H / pdb_00004x9h
Title : Crystal structure of Dscam1 isoform 8.4, N-terminal four Ig domains
Authors : Chen, Q.
Deposited on : 2014-12-11
Resolution : 2.95 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

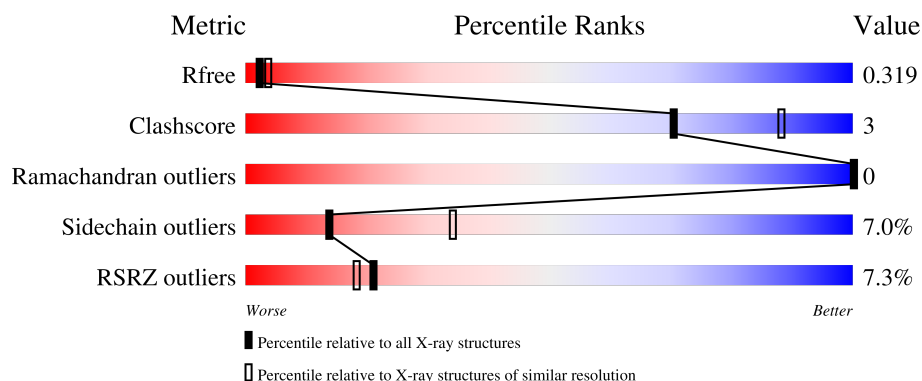
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.95 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	1130 (2.98-2.94)
Clashscore	190562	1157 (2.98-2.94)
Ramachandran outliers	187476	1101 (2.98-2.94)
Sidechain outliers	187428	1101 (2.98-2.94)
RSRZ outliers	180081	1130 (2.98-2.94)

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	400	4% 86% 10% ..
1	B	400	10% 84% 12% ..
2	C	4	25% 75%
2	D	4	50% 50%

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 6231 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 Down syndrome cell adhesion molecule, isoform AP.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	391	Total	C	N	O	S	0	2	0
			3052	1925	530	583	14			
1	B	392	Total	C	N	O	S	0	1	0
			3048	1923	528	583	14			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-3	GLY	-	expression tag	UNP Q0E9L0
A	-2	GLY	-	expression tag	UNP Q0E9L0
B	-3	GLY	-	expression tag	UNP Q0E9L0
B	-2	GLY	-	expression tag	UNP Q0E9L0

- Molecule 2 is an oligosaccharide called alpha-D-mannopyranose-(1-3)-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	C	4	Total	C	N	O	0	0	0
			50	28	2	20			
2	D	4	Total	C	N	O	0	0	0
			50	28	2	20			

- Molecule 3 is 2-acetamido-2-deoxy-beta-D-glucopyranose (CCD ID: NAG) (formula: C₈H₁₅NO₆).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			14	8	1	5		
3	B	1	Total	C	N	O	0	0
			14	8	1	5		

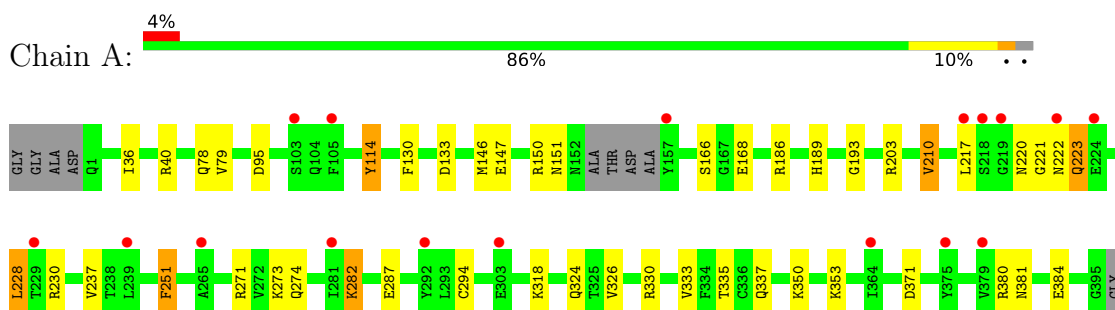
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	1	Total	O	0	0
			1	1		
4	B	2	Total	O	0	0
			2	2		

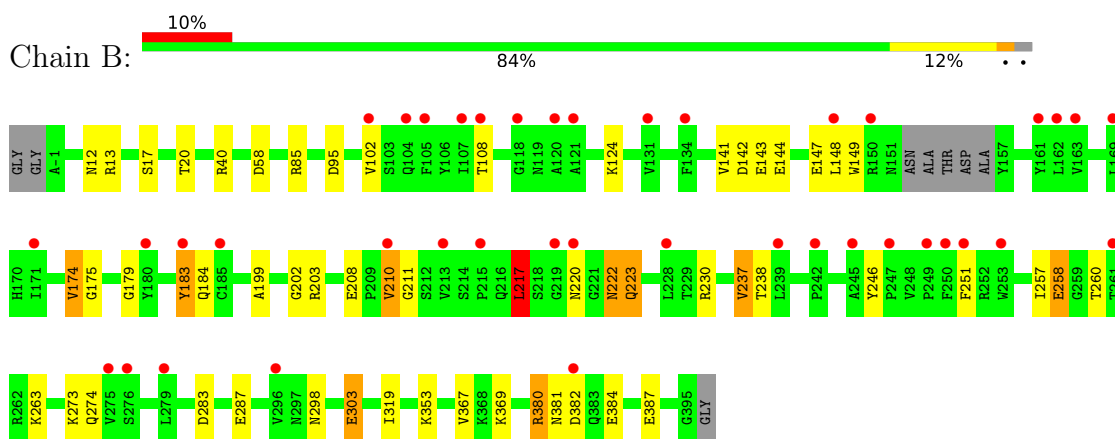
3 Residue-property plots

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: Down syndrome cell adhesion molecule, isoform AP



- Molecule 1: Down syndrome cell adhesion molecule, isoform AP



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



- Molecule 2: alpha-D-mannopyranose-(1-3)-beta-D-mannopyranose-(1-4)-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 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	95.22Å 60.95Å 92.35Å 90.00° 93.56° 90.00°	Depositor
Resolution (Å)	47.52 – 2.95 47.52 – 2.95	Depositor EDS
% Data completeness (in resolution range)	99.3 (47.52-2.95) 99.2 (47.52-2.95)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.93 (at 2.96Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.8.4_1496)	Depositor
R, R_{free}	0.272 , 0.324 0.275 , 0.319	Depositor DCC
R_{free} test set	1149 reflections (5.08%)	wwPDB-VP
Wilson B-factor (Å ²)	80.4	Xtriage
Anisotropy	0.139	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.26 , 0.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.021 for l,-k,h	Xtriage
F_o, F_c correlation	0.90	EDS
Total number of atoms	6231	wwPDB-VP
Average B, all atoms (Å ²)	76.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.17% 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: MAN, NAG, 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	0.30	0/3119	0.65	1/4228 (0.0%)
1	B	0.30	0/3115	0.65	1/4223 (0.0%)
All	All	0.30	0/6234	0.65	2/8451 (0.0%)

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	2

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	A	221	GLY	N-CA-C	7.26	122.39	111.42
1	B	217	LEU	CB-CA-C	-6.60	107.94	115.79

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	220	ASN	Peptide
1	A	223	GLN	Peptide

5.2 Too-close contacts ⓘ

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3052	0	3025	20	0
1	B	3048	0	3022	25	0
2	C	50	0	43	0	0
2	D	50	0	43	0	0
3	A	14	0	13	0	0
3	B	14	0	13	0	0
4	A	1	0	0	0	0
4	B	2	0	0	0	0
All	All	6231	0	6159	42	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 (42) 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:133:ASP:O	1:B:220:ASN:ND2	2.28	0.67
1:A:271:ARG:NH1	1:A:282:LYS:O	2.30	0.65
1:B:246:TYR:O	1:B:298:ASN:ND2	2.32	0.63
1:B:251:PHE:O	1:B:274:GLN:NE2	2.32	0.62
1:B:17:SER:HG	1:B:20:THR:HG1	1.47	0.61
1:B:183:TYR:O	1:B:202:GLY:N	2.35	0.59
1:A:40:ARG:NH1	1:A:78:GLN:OE1	2.39	0.56
1:B:381:ASN:OD1	1:B:382:ASP:N	2.40	0.54
1:A:350:LYS:NZ	1:A:371:ASP:O	2.41	0.53
1:B:95:ASP:OD2	1:B:380:ARG:NH1	2.43	0.52
1:A:287:GLU:N	1:A:287:GLU:OE2	2.41	0.51
1:A:150:ARG:O	1:A:151:ASN:ND2	2.44	0.50
1:A:318:LYS:N	1:A:337:GLN:O	2.45	0.50
1:B:179:GLY:O	1:B:203:ARG:NH1	2.45	0.50
1:B:12:ASN:OD1	1:B:13:ARG:NH2	2.45	0.49
1:B:258:GLU:N	1:B:258:GLU:OE1	2.45	0.49
1:A:114:TYR:N	1:A:114:TYR:CD2	2.79	0.48
1:A:381:ASN:OD1	1:A:384:GLU:N	2.46	0.48
1:B:183:TYR:CD1	1:B:183:TYR:N	2.81	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:222:ASN:OD1	1:A:223:GLN:N	2.40	0.47
1:A:251:PHE:N	1:A:251:PHE:CD1	2.82	0.47
1:A:189:HIS:O	1:A:193:GLY:N	2.47	0.47
1:A:146:MET:HG2	1:A:147:GLU:H	1.81	0.46
1:A:95:ASP:OD2	1:A:380:ARG:NE	2.47	0.45
1:A:324:GLN:NE2	1:A:333:VAL:O	2.48	0.45
1:B:257:ILE:O	1:B:260:THR:OG1	2.22	0.44
1:B:208:GLU:OE1	1:B:208:GLU:N	2.49	0.44
1:B:210:VAL:HG13	1:B:211:GLY:H	1.82	0.44
1:A:222:ASN:CG	1:A:223:GLN:H	2.20	0.44
1:B:283:ASP:N	1:B:283:ASP:OD1	2.50	0.44
1:B:58:ASP:OD1	1:B:58:ASP:N	2.50	0.44
1:A:210:VAL:CG1	1:B:102:VAL:HG11	2.48	0.44
1:B:108:THR:O	1:B:199:ALA:N	2.35	0.42
1:B:141:VAL:HG12	1:B:142:ASP:N	2.35	0.42
1:A:166:SER:OG	1:A:168:GLU:OE1	2.38	0.42
1:B:149:TRP:N	1:B:149:TRP:CD1	2.88	0.41
1:A:228:LEU:O	1:B:223:GLN:HB2	2.21	0.41
1:B:174:VAL:HG13	1:B:175:GLY:N	2.35	0.41
1:B:217:LEU:HD23	1:B:303:GLU:HB2	2.02	0.41
1:A:78:GLN:CG	1:A:79:VAL:N	2.83	0.41
1:B:237:VAL:HG12	1:B:238:THR:H	1.85	0.40
1:B:222:ASN:OD1	1:B:222:ASN:N	2.53	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	389/400 (97%)	370 (95%)	19 (5%)	0	100	100
1	B	389/400 (97%)	366 (94%)	23 (6%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
All	All	778/800 (97%)	736 (95%)	42 (5%)	0	100	100

There are no Ramachandran outliers to report.

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	337/338 (100%)	318 (94%)	19 (6%)	19	43
1	B	336/338 (99%)	308 (92%)	28 (8%)	10	27
All	All	673/676 (100%)	626 (93%)	47 (7%)	14	34

All (47) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	36	ILE
1	A	114	TYR
1	A	130	PHE
1	A	186	ARG
1	A	203	ARG
1	A	210	VAL
1	A	217	LEU
1	A	228	LEU
1	A	230	ARG
1	A	237	VAL
1	A	251	PHE
1	A	273	LYS
1	A	274	GLN
1	A	282	LYS
1	A	294	CYS
1	A	326	VAL
1	A	330	ARG
1	A	335	THR
1	A	353	LYS
1	B	40	ARG

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Mol	Chain	Res	Type
1	B	85	ARG
1	B	124	LYS
1	B	143	GLU
1	B	144	GLU
1	B	147	GLU
1	B	148	LEU
1	B	174	VAL
1	B	183	TYR
1	B	184	GLN
1	B	210	VAL
1	B	217	LEU
1	B	222	ASN
1	B	223	GLN
1	B	230	ARG
1	B	237	VAL
1	B	258	GLU
1	B	263	LYS
1	B	273	LYS
1	B	287	GLU
1	B	303	GLU
1	B	319	ILE
1	B	353	LYS
1	B	367	VAL
1	B	369	LYS
1	B	380	ARG
1	B	384	GLU
1	B	387	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	151	ASN
1	A	152	ASN
1	B	225	HIS

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 ⓘ

8 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	C	1	1,2	14,14,15	0.61	1 (7%)	17,19,21	0.57	0
2	NAG	C	2	2	14,14,15	0.57	0	17,19,21	0.59	0
2	BMA	C	3	2	11,11,12	0.69	0	15,15,17	1.70	4 (26%)
2	MAN	C	4	2	11,11,12	0.69	1 (9%)	15,15,17	1.58	3 (20%)
2	NAG	D	1	1,2	14,14,15	0.29	0	17,19,21	0.47	0
2	NAG	D	2	2	14,14,15	0.27	0	17,19,21	0.41	0
2	BMA	D	3	2	11,11,12	0.81	0	15,15,17	1.18	2 (13%)
2	MAN	D	4	2	11,11,12	1.00	1 (9%)	15,15,17	1.72	3 (20%)

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

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	4	MAN	O5-C5	2.52	1.48	1.43
2	C	1	NAG	O5-C1	-2.22	1.40	1.43
2	C	4	MAN	C1-C2	2.09	1.57	1.52

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	4	MAN	C1-O5-C5	4.82	118.65	112.19
2	C	4	MAN	C1-O5-C5	4.26	117.90	112.19
2	C	3	BMA	C1-O5-C5	4.25	117.89	112.19
2	D	3	BMA	C1-O5-C5	2.69	115.79	112.19
2	C	3	BMA	O5-C1-C2	2.61	117.02	110.79
2	C	4	MAN	O5-C1-C2	2.44	116.60	110.79
2	D	3	BMA	C1-C2-C3	-2.42	106.12	109.64
2	D	4	MAN	O2-C2-C3	-2.41	105.15	110.15
2	C	3	BMA	O2-C2-C3	-2.31	105.37	110.15
2	D	4	MAN	O5-C1-C2	2.26	116.19	110.79
2	C	4	MAN	O2-C2-C3	-2.23	105.53	110.15
2	C	3	BMA	C1-C2-C3	2.22	112.87	109.64

There are no chirality outliers.

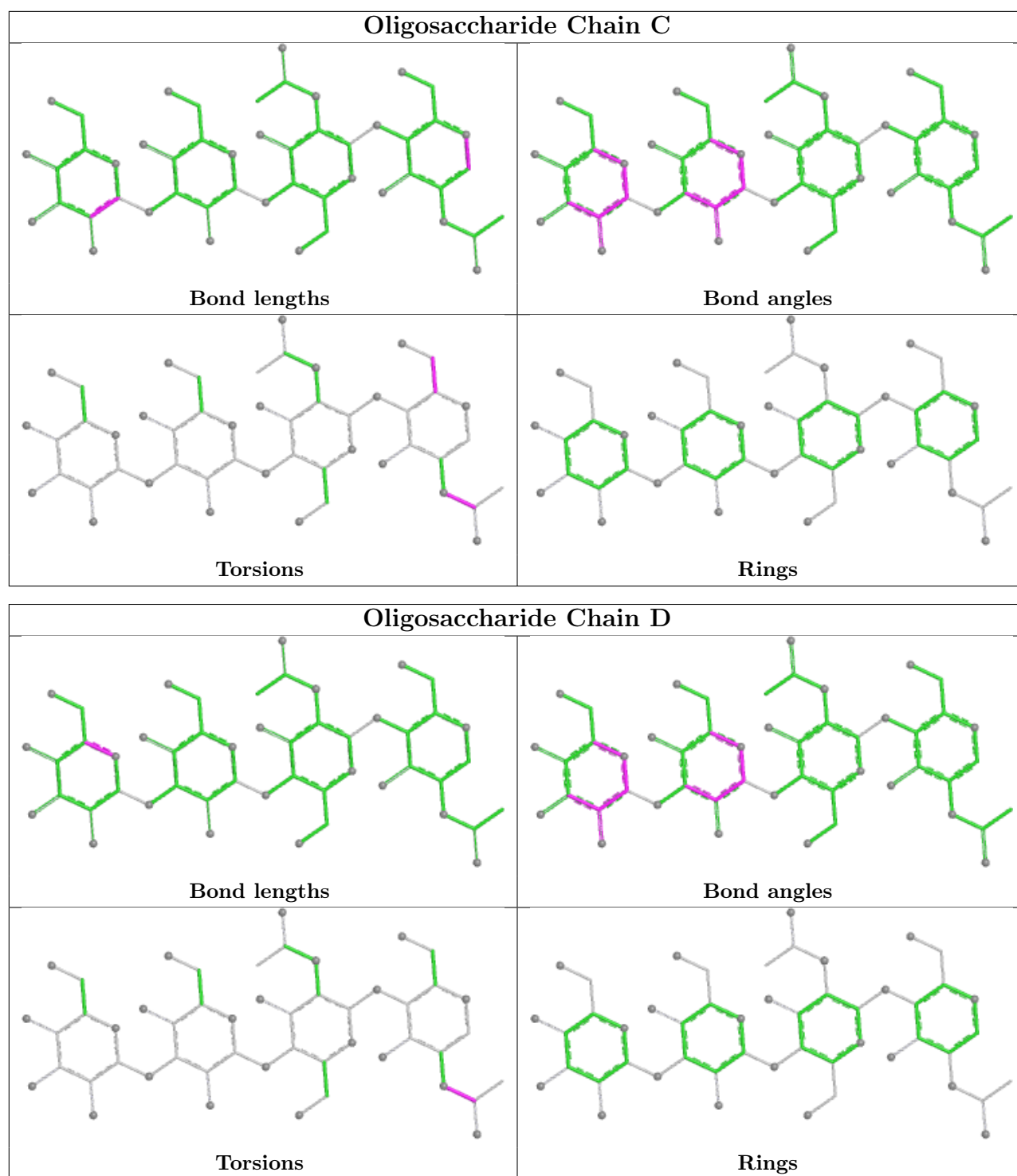
All (6) torsion outliers are listed below:

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

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 [i](#)

2 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The

Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
3	NAG	B	405	1	14,14,15	0.29	0	17,19,21	0.43	0
3	NAG	A	405	1	14,14,15	0.26	0	17,19,21	0.39	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	NAG	B	405	1	-	0/6/23/26	0/1/1/1
3	NAG	A	405	1	-	1/6/23/26	0/1/1/1

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (1) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	405	NAG	O5-C5-C6-O6

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	391/400 (97%)	0.47	17 (4%)	40	32	28, 71, 97, 118	2 (0%)
1	B	392/400 (98%)	0.80	40 (10%)	12	11	50, 77, 111, 150	1 (0%)
All	All	783/800 (97%)	0.63	57 (7%)	21	18	28, 73, 108, 150	3 (0%)

All (57) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	161	TYR	4.6
1	B	276	SER	4.3
1	A	222	ASN	4.0
1	B	253	TRP	4.0
1	B	220	ASN	3.9
1	B	245	ALA	3.9
1	B	162	LEU	3.7
1	B	219	GLY	3.7
1	A	103	SER	3.7
1	B	108	THR	3.6
1	B	163	VAL	3.6
1	B	171	ILE	3.2
1	A	217	LEU	3.2
1	A	281	ILE	3.2
1	B	251	PHE	3.1
1	B	210	VAL	3.0
1	B	105	PHE	2.9
1	B	107	ILE	2.9
1	B	169	LEU	2.9
1	B	183	TYR	2.8
1	A	375	TYR	2.8
1	B	249	PRO	2.7
1	B	239	LEU	2.6
1	B	148	LEU	2.6

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Mol	Chain	Res	Type	RSRZ
1	B	215	PRO	2.5
1	B	242	PRO	2.5
1	B	134	PHE	2.5
1	B	120	ALA	2.5
1	B	118	GLY	2.5
1	B	213	VAL	2.5
1	A	364	ILE	2.4
1	A	105	PHE	2.4
1	B	296	VAL	2.4
1	B	121	ALA	2.4
1	A	157	TYR	2.4
1	B	185	CYS	2.3
1	A	239	LEU	2.3
1	A	292	TYR	2.2
1	B	180	TYR	2.2
1	B	228	LEU	2.2
1	B	104	GLN	2.2
1	A	379	VAL	2.2
1	B	102	VAL	2.1
1	B	250	PHE	2.1
1	A	218	SER	2.1
1	B	131	VAL	2.1
1	B	382	ASP	2.1
1	A	265	ALA	2.1
1	A	224	GLU	2.1
1	A	219	GLY	2.0
1	A	229	THR	2.0
1	B	261	THR	2.0
1	A	303	GLU	2.0
1	B	275	VAL	2.0
1	B	247	PRO	2.0
1	B	279	LEU	2.0
1	B	150	ARG	2.0

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

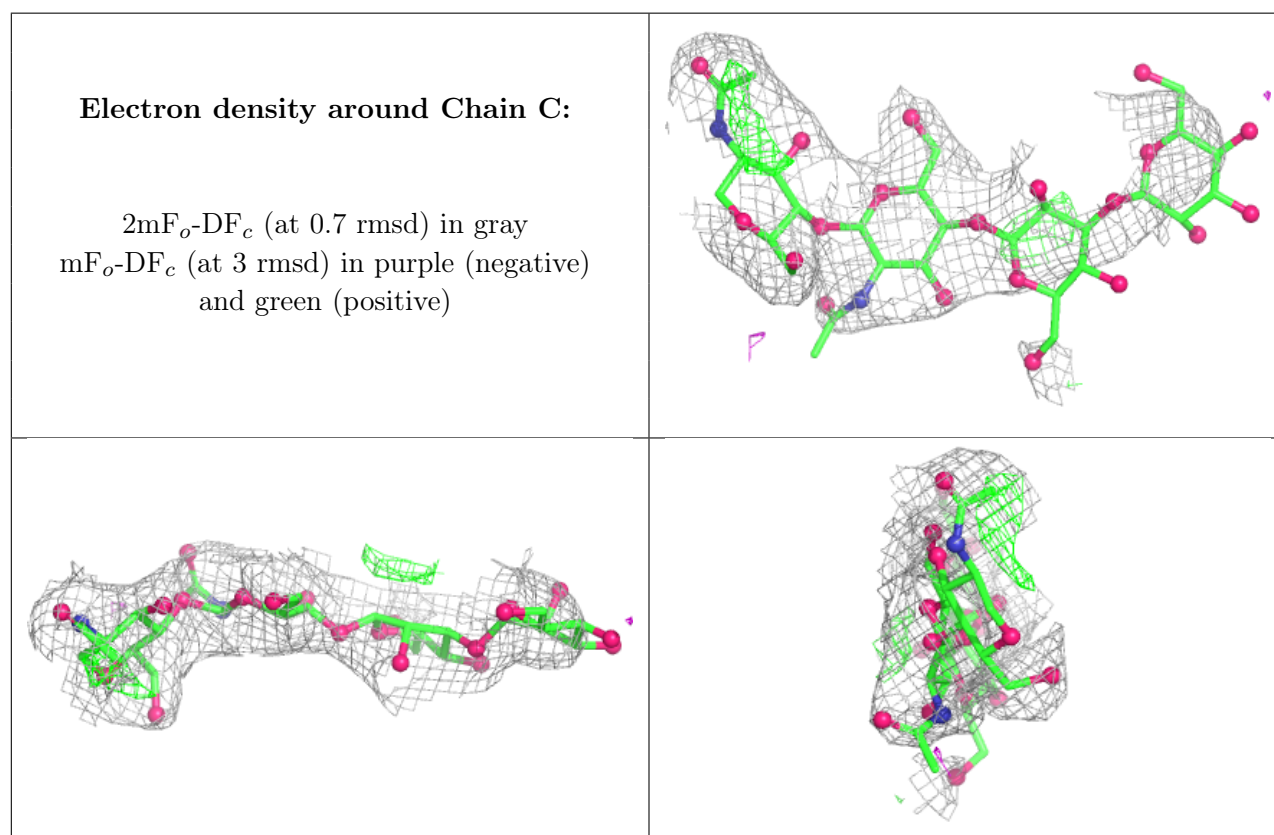
There are no non-standard protein/DNA/RNA residues in this entry.

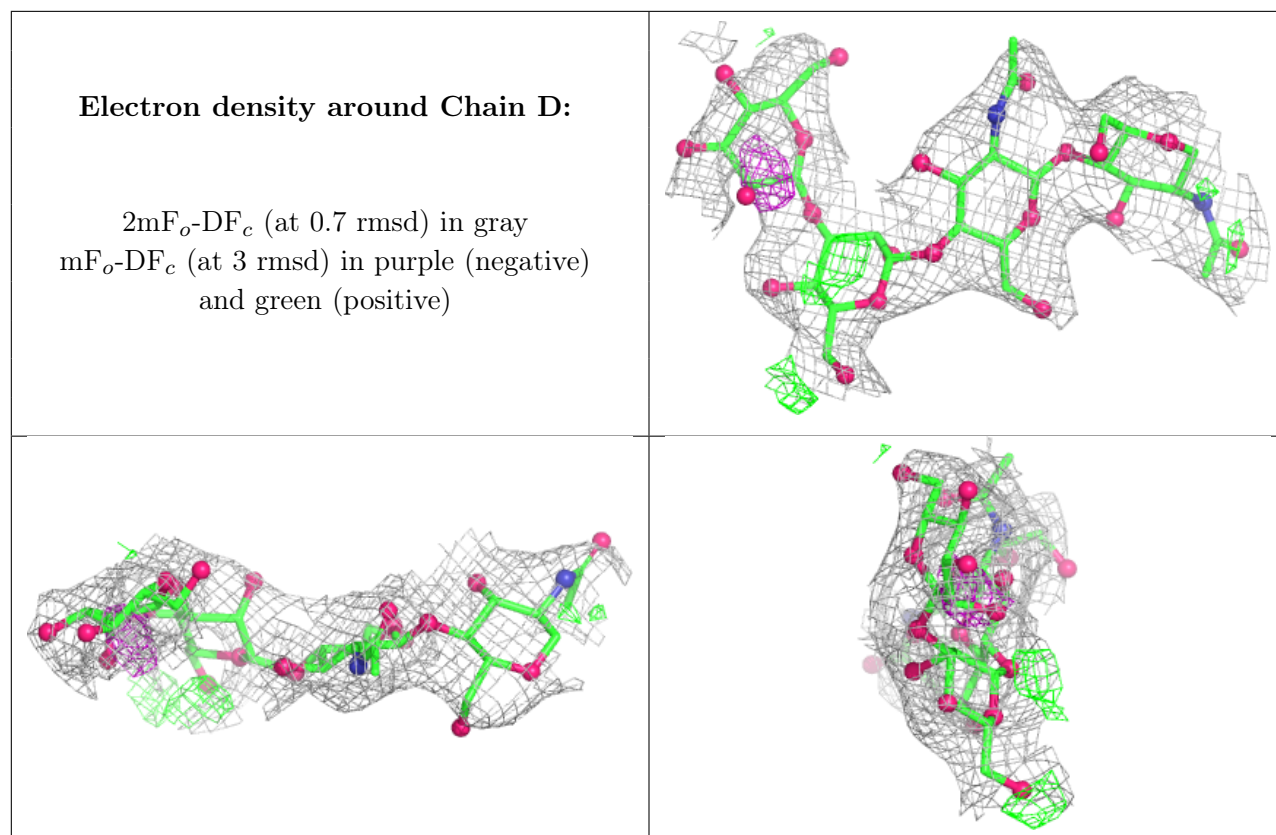
6.3 Carbohydrates [i](#)

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	MAN	C	4	11/12	0.38	0.13	115,137,142,142	0
2	BMA	C	3	11/12	0.40	0.15	120,128,137,144	0
2	MAN	D	4	11/12	0.49	0.20	78,84,91,93	0
2	BMA	D	3	11/12	0.63	0.14	73,80,86,87	0
2	NAG	C	1	14/15	0.77	0.13	54,66,74,76	0
2	NAG	D	1	14/15	0.77	0.14	63,72,77,83	0
2	NAG	C	2	14/15	0.80	0.12	81,93,106,114	0
2	NAG	D	2	14/15	0.87	0.10	73,78,82,85	0

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





6.4 Ligands [i](#)

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	NAG	B	405	14/15	0.49	0.14	106,117,121,121	0
3	NAG	A	405	14/15	0.67	0.15	83,88,92,94	0

6.5 Other polymers [i](#)

There are no such residues in this entry.