



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 10, 2024 – 08:52 AM EDT

PDB ID : 4RM0
Title : Crystal structure of Norovirus OIF P domain in complex with Lewis a trisaccharide
Authors : Liu, W.; Chen, Y.; Tan, M.; Xia, M.; Li, X.; Jiang, X.; Rao, Z.
Deposited on : 2014-10-18
Resolution : 2.00 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 : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.36
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.36

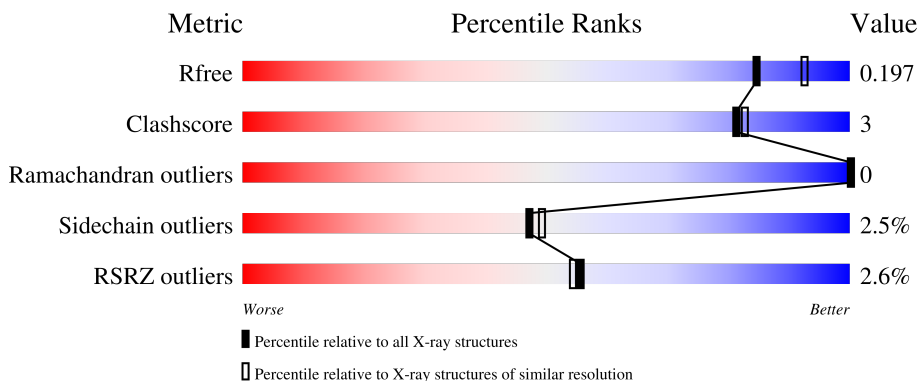
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.00 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}	130704	8085 (2.00-2.00)
Clashscore	141614	9178 (2.00-2.00)
Ramachandran outliers	138981	9054 (2.00-2.00)
Sidechain outliers	138945	9053 (2.00-2.00)
RSRZ outliers	127900	7900 (2.00-2.00)

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	316	 91% 6%
1	B	316	 88% 9%
2	C	3	 33% 67%
3	D	3	 100%

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 5346 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 Capsid protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	307	2371	1506	406	454	5	0	0	0
1	B	306	2362	1500	404	453	5	0	0	0

There are 30 discrepancies between the modelled and reference sequences:

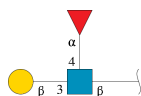
Chain	Residue	Modelled	Actual	Comment	Reference
A	212	GLY	-	expression tag	UNP Q6B7R3
A	213	PRO	-	expression tag	UNP Q6B7R3
A	214	LEU	-	expression tag	UNP Q6B7R3
A	215	GLY	-	expression tag	UNP Q6B7R3
A	216	SER	-	expression tag	UNP Q6B7R3
A	217	PRO	-	expression tag	UNP Q6B7R3
A	218	GLU	-	expression tag	UNP Q6B7R3
A	219	PHE	-	expression tag	UNP Q6B7R3
A	375	ASN	HIS	engineered mutation	UNP Q6B7R3
A	376	ILE	HIS	engineered mutation	UNP Q6B7R3
A	377	ALA	SER	engineered mutation	UNP Q6B7R3
A	378	SER	GLN	engineered mutation	UNP Q6B7R3
A	379	ASN	HIS	engineered mutation	UNP Q6B7R3
A	387	VAL	LEU	engineered mutation	UNP Q6B7R3
A	389	ILE	VAL	engineered mutation	UNP Q6B7R3
B	212	GLY	-	expression tag	UNP Q6B7R3
B	213	PRO	-	expression tag	UNP Q6B7R3
B	214	LEU	-	expression tag	UNP Q6B7R3
B	215	GLY	-	expression tag	UNP Q6B7R3
B	216	SER	-	expression tag	UNP Q6B7R3
B	217	PRO	-	expression tag	UNP Q6B7R3
B	218	GLU	-	expression tag	UNP Q6B7R3
B	219	PHE	-	expression tag	UNP Q6B7R3
B	375	ASN	HIS	engineered mutation	UNP Q6B7R3
B	376	ILE	HIS	engineered mutation	UNP Q6B7R3

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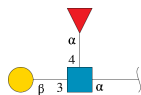
Chain	Residue	Modelled	Actual	Comment	Reference
B	377	ALA	SER	engineered mutation	UNP Q6B7R3
B	378	SER	GLN	engineered mutation	UNP Q6B7R3
B	379	ASN	HIS	engineered mutation	UNP Q6B7R3
B	387	VAL	LEU	engineered mutation	UNP Q6B7R3
B	389	ILE	VAL	engineered mutation	UNP Q6B7R3

- Molecule 2 is an oligosaccharide called beta-D-galactopyranose-(1-3)-[alpha-L-fucopyranose-(1-4)]2-acetamido-2-deoxy-beta-D-glucopyranose.



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
			Total	C	N	O			
2	C	3	36	20	1	15	0	0	0

- Molecule 3 is an oligosaccharide called beta-D-galactopyranose-(1-3)-[alpha-L-fucopyranose-(1-4)]2-acetamido-2-deoxy-alpha-D-glucopyranose.



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
			Total	C	N	O			
3	D	3	36	20	1	15	0	0	0

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	282	Total	O	0	0
			282	282		
4	B	259	Total	O	0	0
			259	259		

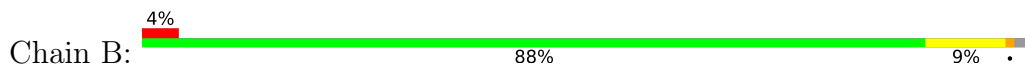
3 Residue-property plots [i](#)

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: Capsid protein



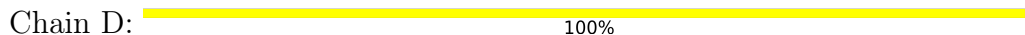
- Molecule 1: Capsid protein



- Molecule 2: beta-D-galactopyranose-(1-3)-[alpha-L-fucopyranose-(1-4)]2-acetamido-2-deoxy-beta-D-glucopyranose



- Molecule 3: beta-D-galactopyranose-(1-3)-[alpha-L-fucopyranose-(1-4)]2-acetamido-2-deoxy-alpha-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	54.85Å 84.16Å 124.74Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	25.05 – 2.00 25.05 – 2.00	Depositor EDS
% Data completeness (in resolution range)	94.7 (25.05-2.00) 94.7 (25.05-2.00)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.67 (at 1.99Å)	Xtrriage
Refinement program	PHENIX (phenix.refine: 1.7_650)	Depositor
R, R_{free}	0.171 , 0.209 0.162 , 0.197	Depositor DCC
R_{free} test set	1884 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	21.0	Xtrriage
Anisotropy	0.175	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 47.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	5346	wwPDB-VP
Average B, all atoms (Å ²)	21.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.27% 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: NDG, FUC, GAL, NAG

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
1	A	0.41	0/2439	0.53	0/3342
1	B	0.40	0/2430	0.53	0/3331
All	All	0.41	0/4869	0.53	0/6673

There are no bond length outliers.

There are no bond angle outliers.

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	2371	0	2287	7	0
1	B	2362	0	2274	16	0
2	C	36	0	33	1	0
3	D	36	0	30	0	0
4	A	282	0	0	1	0
4	B	259	0	0	1	0
All	All	5346	0	4624	24	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 (24) 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:B:407:GLY:O	1:B:410:THR:HG22	1.65	0.97
1:B:406:SER:HB2	1:B:410:THR:HG21	1.49	0.92
1:B:406:SER:CB	1:B:410:THR:HG21	2.24	0.68
1:B:406:SER:HB2	1:B:410:THR:CG2	2.23	0.64
1:A:502:ASN:ND2	4:A:837:HOH:O	2.30	0.64
1:B:266:LEU:HG	1:B:490:ARG:HA	1.83	0.60
1:B:485:GLU:HG3	1:B:524:LEU:HD22	1.87	0.56
2:C:1:NAG:H61	2:C:3:FUC:C1	2.39	0.52
1:B:520:GLN:HG2	1:B:521:PHE:CE1	2.46	0.51
1:B:252:ASN:HB3	1:B:255:VAL:HG22	1.93	0.51
1:B:466:GLN:HB2	1:B:517:TRP:CG	2.48	0.49
1:B:487:LYS:HE3	1:B:522:TYR:O	2.13	0.48
1:B:473:ARG:HD2	1:B:515:ASP:HB2	1.96	0.46
1:A:403:PRO:HG2	1:A:405:TYR:CE1	2.50	0.46
1:A:237:ASN:HB3	1:A:240:PHE:O	2.15	0.46
1:A:420:GLY:HA2	1:A:428:ILE:HD11	1.97	0.46
1:B:355:SER:HB2	1:B:403:PRO:HB3	1.99	0.45
1:B:223:LYS:HA	1:B:224:PRO:HD3	1.84	0.45
1:B:410:THR:HG23	4:B:728:HOH:O	2.17	0.44
1:B:297:ASP:HB2	1:B:359:LYS:HD2	2.00	0.43
1:A:312:ASP:OD2	1:A:314:THR:OG1	2.26	0.43
1:B:237:ASN:HB3	1:B:240:PHE:O	2.19	0.42
1:A:507:PRO:HA	1:A:508:PRO:HD3	1.96	0.41
1:A:473:ARG:HG2	1:A:485:GLU:HG2	2.02	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	305/316 (96%)	290 (95%)	15 (5%)	0	100 100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	B	304/316 (96%)	291 (96%)	13 (4%)	0	100	100
All	All	609/632 (96%)	581 (95%)	28 (5%)	0	100	100

There are no Ramachandran outliers to report.

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	261/268 (97%)	254 (97%)	7 (3%)	44	46
1	B	260/268 (97%)	254 (98%)	6 (2%)	50	53
All	All	521/536 (97%)	508 (98%)	13 (2%)	47	49

All (13) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	222	THR
1	A	223	LYS
1	A	293	ARG
1	A	312	ASP
1	A	411	LEU
1	A	490	ARG
1	A	527	MET
1	B	293	ARG
1	B	341	SER
1	B	381	GLN
1	B	473	ARG
1	B	520	GLN
1	B	527	MET

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	502	ASN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

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	NAG	C	1	2	15,15,15	2.27	5 (33%)	21,21,21	1.30	1 (4%)
2	GAL	C	2	2	11,11,12	1.72	3 (27%)	15,15,17	1.38	1 (6%)
2	FUC	C	3	2	10,10,11	2.10	3 (30%)	14,14,16	1.06	2 (14%)
3	NDG	D	1	3	15,15,15	2.23	5 (33%)	21,21,21	1.29	2 (9%)
3	GAL	D	2	3	11,11,12	1.74	3 (27%)	15,15,17	1.16	1 (6%)
3	FUC	D	3	3	10,10,11	2.29	5 (50%)	14,14,16	1.27	2 (14%)

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	2	-	1/6/26/26	0/1/1/1
2	GAL	C	2	2	-	0/2/19/22	0/1/1/1
2	FUC	C	3	2	-	-	0/1/1/1
3	NDG	D	1	3	-	0/6/26/26	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	GAL	D	2	3	-	0/2/19/22	0/1/1/1
3	FUC	D	3	3	-	-	0/1/1/1

All (24) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	1	NAG	C7-N2	4.50	1.49	1.34
3	D	3	FUC	C2-C3	-4.32	1.46	1.52
3	D	1	NDG	C7-N2	4.26	1.49	1.34
3	D	3	FUC	C4-C3	-4.08	1.41	1.52
2	C	1	NAG	C4-C3	-3.94	1.42	1.52
2	C	3	FUC	C4-C3	-3.92	1.42	1.52
3	D	1	NDG	C4-C3	-3.91	1.42	1.52
3	D	1	NDG	O3-C3	3.64	1.51	1.43
2	C	3	FUC	C2-C3	-3.59	1.47	1.52
2	C	1	NAG	O3-C3	3.44	1.51	1.43
3	D	1	NDG	C3-C2	-3.24	1.47	1.53
2	C	1	NAG	C3-C2	-3.02	1.47	1.53
3	D	2	GAL	C2-C3	-2.97	1.48	1.52
2	C	2	GAL	C4-C3	-2.85	1.45	1.52
2	C	2	GAL	C2-C3	-2.79	1.48	1.52
2	C	1	NAG	C2-N2	2.71	1.50	1.45
3	D	2	GAL	O5-C5	2.61	1.48	1.43
3	D	2	GAL	C4-C3	-2.56	1.45	1.52
3	D	1	NDG	C2-N2	2.34	1.49	1.45
3	D	3	FUC	C6-C5	-2.31	1.46	1.51
2	C	2	GAL	O5-C5	2.31	1.48	1.43
2	C	3	FUC	C6-C5	-2.25	1.46	1.51
3	D	3	FUC	O5-C1	-2.10	1.40	1.43
3	D	3	FUC	O5-C5	2.01	1.47	1.43

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	2	GAL	C1-C2-C3	3.89	114.44	109.67
3	D	2	GAL	C1-C2-C3	3.00	113.35	109.67
2	C	1	NAG	O5-C1-C2	2.85	112.38	109.52
3	D	3	FUC	C3-C4-C5	2.48	113.64	109.77
3	D	1	NDG	O5-C1-C2	2.38	111.91	109.52
2	C	3	FUC	C3-C4-C5	2.30	113.36	109.77
3	D	1	NDG	C8-C7-N2	2.20	119.83	116.10
3	D	3	FUC	O5-C5-C4	2.17	113.42	109.52

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	3	FUC	O5-C5-C4	2.05	113.20	109.52

There are no chirality outliers.

All (1) torsion outliers are listed below:

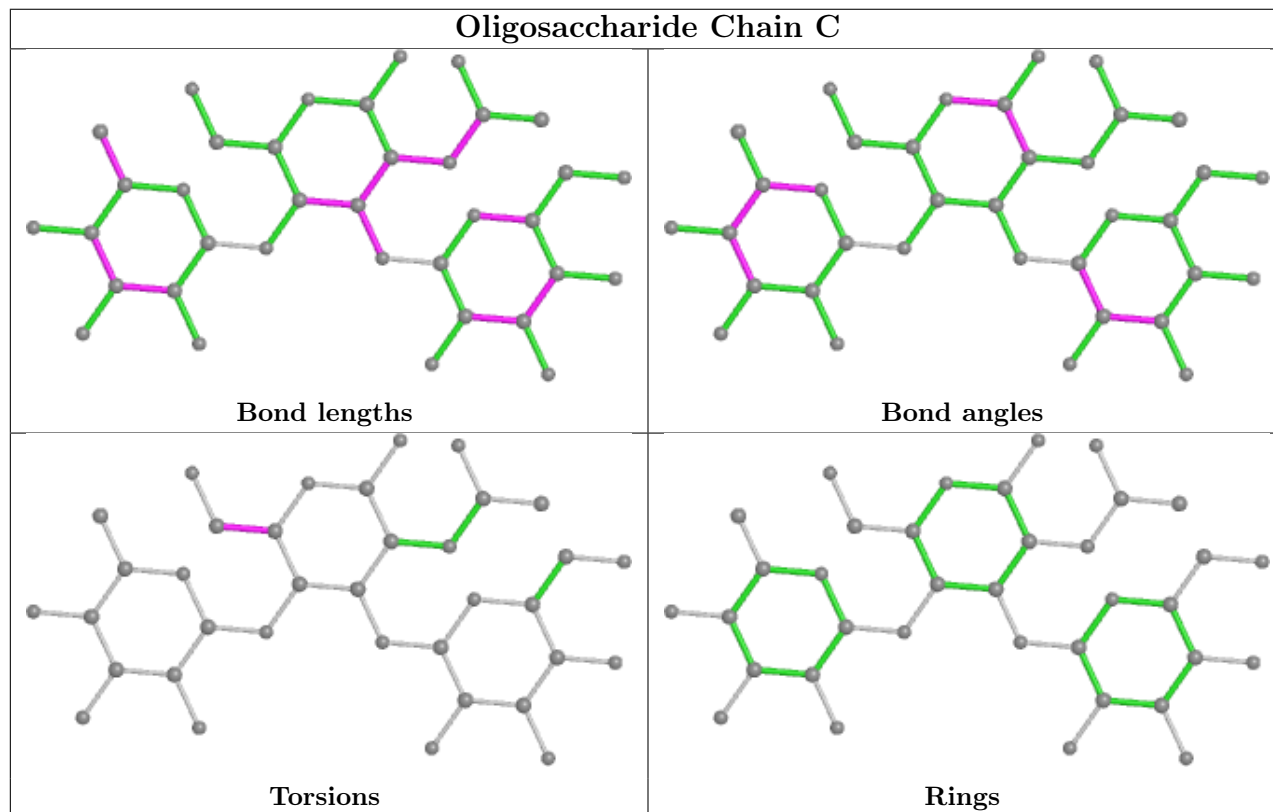
Mol	Chain	Res	Type	Atoms
2	C	1	NAG	C4-C5-C6-O6

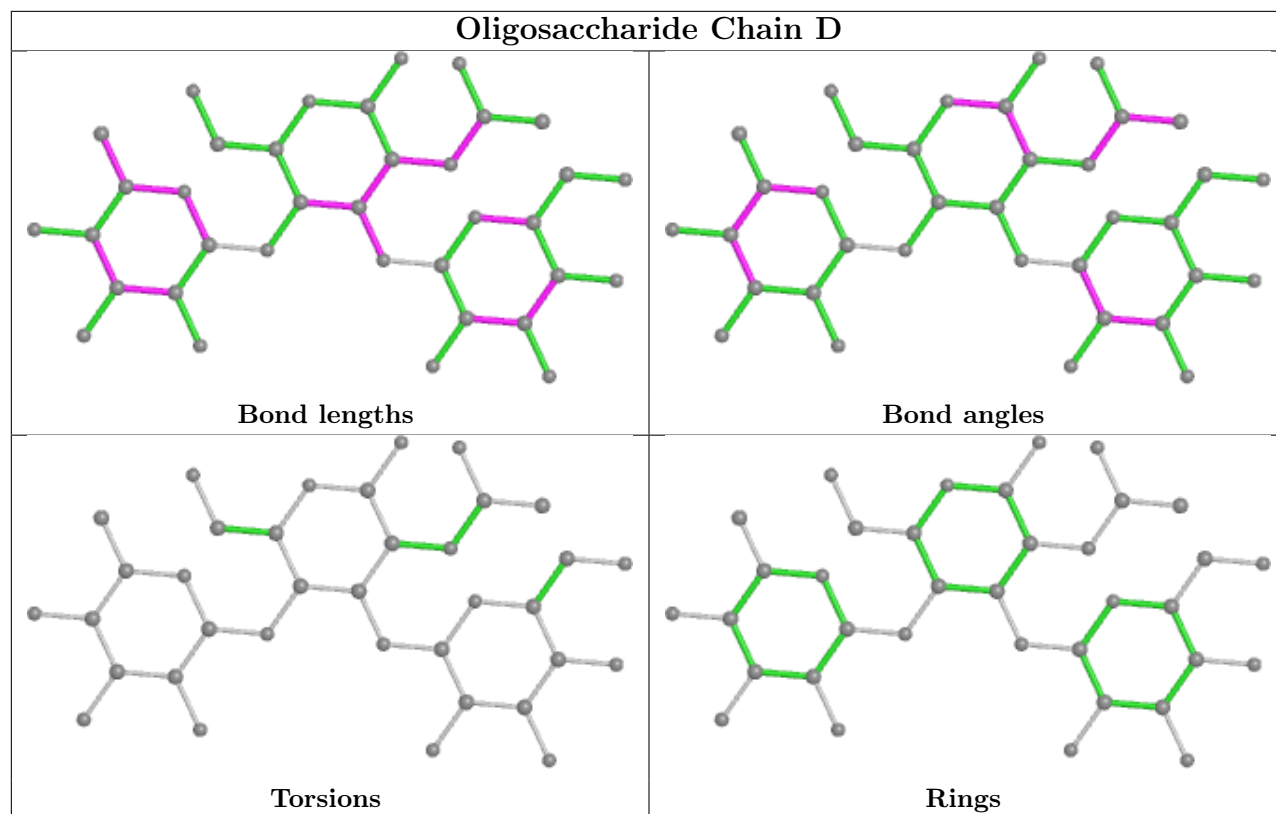
There are no ring outliers.

2 monomers are involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	1	NAG	1	0
2	C	3	FUC	1	0

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





5.6 Ligand geometry [i](#)

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	307/316 (97%)	0.09	4 (1%) 77 76	12, 18, 30, 44	0
1	B	306/316 (96%)	0.28	12 (3%) 39 38	13, 19, 34, 47	0
All	All	613/632 (96%)	0.18	16 (2%) 56 54	12, 19, 33, 47	0

All (16) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	295	TYR	6.8
1	B	254	ASP	5.3
1	A	253	ALA	4.5
1	B	222	THR	4.4
1	B	527	MET	3.8
1	B	521	PHE	3.6
1	A	222	THR	3.3
1	A	494	ILE	3.1
1	B	342	GLY	3.0
1	B	255	VAL	3.0
1	B	502	ASN	2.4
1	B	341	SER	2.3
1	B	343	ASP	2.3
1	B	494	ILE	2.2
1	B	525	ALA	2.1
1	B	473	ARG	2.0

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

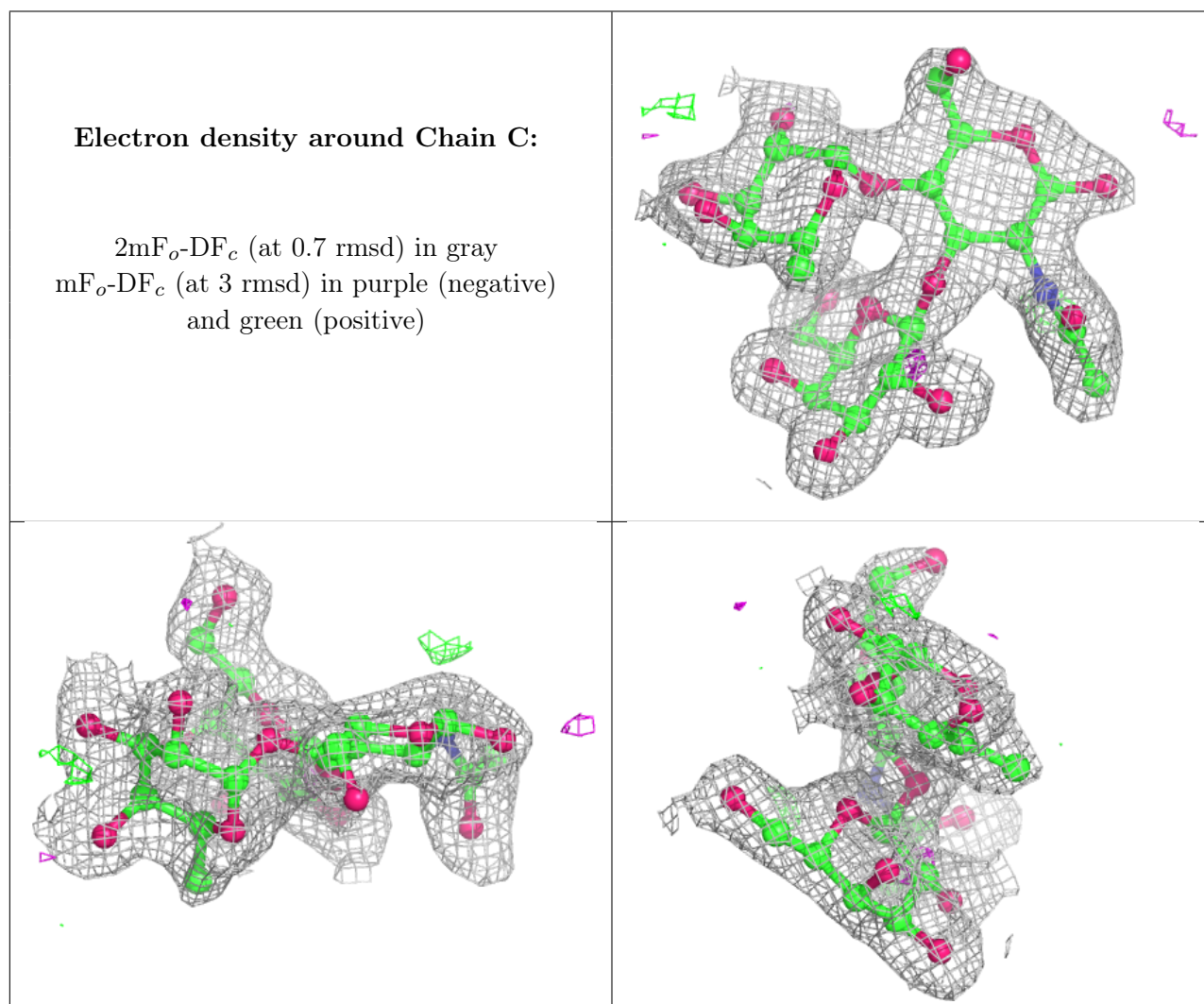
There are no non-standard protein/DNA/RNA residues in this entry.

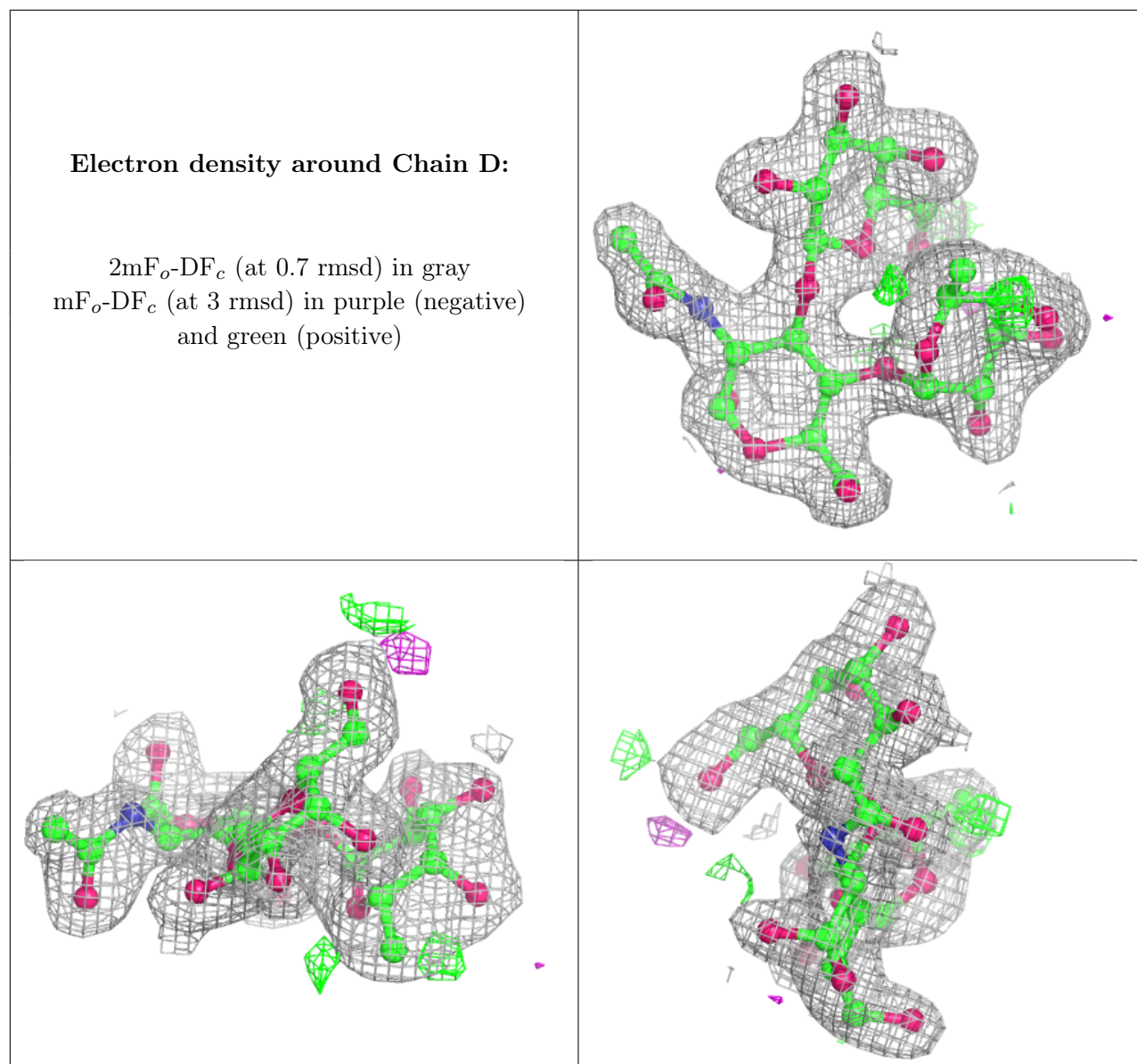
6.3 Carbohydrates [i](#)

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
3	FUC	D	3	10/11	0.90	0.15	22,26,35,36	0
2	GAL	C	2	11/12	0.94	0.12	16,21,25,27	0
2	FUC	C	3	10/11	0.94	0.15	27,30,33,35	0
2	NAG	C	1	15/15	0.94	0.14	25,31,34,35	4
3	GAL	D	2	11/12	0.95	0.09	12,16,18,18	0
3	NDG	D	1	15/15	0.95	0.13	19,22,24,24	3

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](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.