



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 08:13 AM UTC

PDB ID : 7JPI / pdb_00007jpi
Title : Crystal structure of EBOV glycoprotein with modified HR2 stalk at 2.3Å resolution
Authors : Chaudhary, A.; Stanfield, R.L.; Wilson, I.A.; Zhu, J.
Deposited on : 2020-08-08
Resolution : 2.28 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

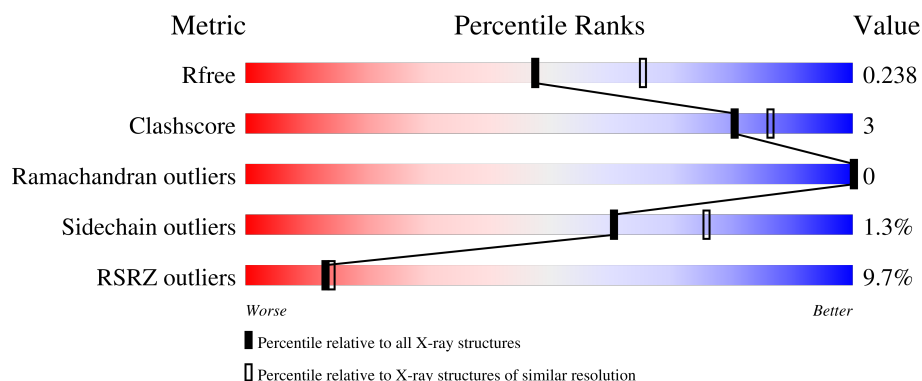
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.28 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	9078 (2.30-2.26)
Clashscore	190562	9802 (2.30-2.26)
Ramachandran outliers	187476	9690 (2.30-2.26)
Sidechain outliers	187428	9691 (2.30-2.26)
RSRZ outliers	180081	9085 (2.30-2.26)

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	297	9% 76% 7% 16%
2	B	165	5% 72% . 24%
3	F	2	100%
3	H	2	100%
4	C	6	50% 50%

2 Entry composition

There are 8 unique types of molecules in this entry. The entry contains 3239 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 Envelope glycoprotein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	249	Total	C	N	O	S	0	1	0
			1960	1254	331	370	5			

- Molecule 2 is a protein called Envelope glycoprotein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	125	Total	C	N	O	S	0	0	0
			982	624	170	182	6			

There are 30 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	615	LEU	TRP	engineered mutation	UNP Q05320
B	638	ALA	-	expression tag	UNP Q05320
B	639	SER	-	expression tag	UNP Q05320
B	640	GLY	-	expression tag	UNP Q05320
B	641	TYR	-	expression tag	UNP Q05320
B	642	ILE	-	expression tag	UNP Q05320
B	643	PRO	-	expression tag	UNP Q05320
B	644	GLU	-	expression tag	UNP Q05320
B	645	ALA	-	expression tag	UNP Q05320
B	646	PRO	-	expression tag	UNP Q05320
B	647	ARG	-	expression tag	UNP Q05320
B	648	ASP	-	expression tag	UNP Q05320
B	649	GLY	-	expression tag	UNP Q05320
B	650	GLN	-	expression tag	UNP Q05320
B	651	ALA	-	expression tag	UNP Q05320
B	652	TYR	-	expression tag	UNP Q05320
B	653	VAL	-	expression tag	UNP Q05320
B	654	ARG	-	expression tag	UNP Q05320
B	655	LYS	-	expression tag	UNP Q05320
B	656	ASP	-	expression tag	UNP Q05320

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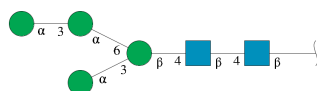
Chain	Residue	Modelled	Actual	Comment	Reference
B	657	GLY	-	expression tag	UNP Q05320
B	658	GLU	-	expression tag	UNP Q05320
B	659	TRP	-	expression tag	UNP Q05320
B	660	VAL	-	expression tag	UNP Q05320
B	661	LEU	-	expression tag	UNP Q05320
B	662	LEU	-	expression tag	UNP Q05320
B	663	SER	-	expression tag	UNP Q05320
B	664	THR	-	expression tag	UNP Q05320
B	665	PHE	-	expression tag	UNP Q05320
B	666	LEU	-	expression tag	UNP Q05320

- 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	F	2	Total	C	N	O	0	0	0
			28	16	2	10			
3	H	2	Total	C	N	O	0	0	0
			28	16	2	10			

- Molecule 4 is an oligosaccharide called alpha-D-mannopyranose-(1-3)-alpha-D-mannopyranose-(1-6)-[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
4	C	6	Total	C	N	O	0	0	0
			72	40	2	30			

- Molecule 5 is GLYCEROL (CCD ID: GOL) (formula: C₃H₈O₃).



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

- Molecule 6 is 2-acetamido-2-deoxy-beta-D-glucopyranose (CCD ID: NAG) (formula: $C_8H_{15}NO_6$).



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

- Molecule 7 is PHOSPHATE ION (CCD ID: PO4) (formula: O_4P).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
7	A	1	Total	O	P	0	0
			5	4	1		

- Molecule 8 is water.

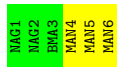
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	A	90	Total	O	0	0
			90	90		
8	B	48	Total	O	0	0
			48	48		

ido-2-deoxy-beta-D-glucopyranose

Chain C:

50%

50%



4 Data and refinement statistics

Property	Value	Source
Space group	H 3 2	Depositor
Cell constants a, b, c, α , β , γ	114.58Å 114.58Å 312.38Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	49.00 – 2.28 49.00 – 2.28	Depositor EDS
% Data completeness (in resolution range)	95.7 (49.00-2.28) 95.7 (49.00-2.28)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	0.89 (at 2.27Å)	Xtriage
Refinement program	PHENIX 1.16_3549	Depositor
R, R_{free}	0.200 , 0.237 0.201 , 0.238	Depositor DCC
R_{free} test set	1715 reflections (4.71%)	wwPDB-VP
Wilson B-factor (Å ²)	42.9	Xtriage
Anisotropy	0.350	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 46.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	3239	wwPDB-VP
Average B, all atoms (Å ²)	58.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.32% 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, PO4, NAG, 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	0.13	0/1993	0.32	0/2707
2	B	0.11	0/1004	0.30	0/1367
All	All	0.12	0/2997	0.32	0/4074

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	1960	0	1879	14	0
2	B	982	0	960	4	0
3	F	28	0	25	1	0
3	H	28	0	25	0	0
4	C	72	0	61	0	0
5	A	12	0	16	0	0
6	A	14	0	13	0	0
7	A	5	0	0	0	0
8	A	90	0	0	0	0
8	B	48	0	0	0	0
All	All	3239	0	2979	16	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 (16) 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:261:TYR:HA	1:A:266:ARG:HD2	1.83	0.60
1:A:44:GLN:HA	2:B:504:ILE:HG13	1.84	0.59
1:A:106:GLU:HG3	1:A:290:PHE:CD1	2.38	0.58
1:A:50:LYS:HZ3	2:B:502:GLU:N	2.03	0.56
2:B:618:ASN:O	2:B:622:LYS:HG2	2.09	0.53
1:A:61:ASN:O	1:A:64:ARG:NH2	2.45	0.49
1:A:35:LEU:HB3	1:A:46:SER:O	2.15	0.46
1:A:229:GLU:HG3	1:A:229:GLU:O	2.17	0.44
1:A:267:SER:HB2	1:A:273:LEU:HB3	1.99	0.43
1:A:46:SER:HA	1:A:50:LYS:HD3	1.99	0.43
1:A:64:ARG:NH1	1:A:193:PHE:O	2.52	0.43
1:A:260:ILE:HG23	1:A:265:LYS:HB2	2.00	0.43
1:A:56:LYS:HD2	1:A:188:GLN:NE2	2.35	0.42
1:A:130:ARG:O	1:A:162:TYR:HB3	2.19	0.42
1:A:194:PHE:HE1	2:B:515:LEU:HD23	1.86	0.41
3:F:1:NAG:H62	3:F:2:NAG:C7	2.52	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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	240/297 (81%)	233 (97%)	7 (3%)	0	100	100
2	B	123/165 (74%)	119 (97%)	4 (3%)	0	100	100
All	All	363/462 (79%)	352 (97%)	11 (3%)	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	209/252 (83%)	207 (99%)	2 (1%)	68	81
2	B	104/138 (75%)	102 (98%)	2 (2%)	50	66
All	All	313/390 (80%)	309 (99%)	4 (1%)	61	76

All (4) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	64	ARG
1	A	70	LEU
2	B	623	ILE
2	B	625	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

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](#)

10 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	$\# Z > 2$
4	NAG	C	1	4,2	14,14,15	0.20	0	17,19,21	0.53	0
4	NAG	C	2	4	14,14,15	0.16	0	17,19,21	0.55	0
4	BMA	C	3	4	11,11,12	0.45	0	15,15,17	0.74	0
4	MAN	C	4	4	11,11,12	0.73	0	15,15,17	0.94	1 (6%)
4	MAN	C	5	4	11,11,12	0.62	0	15,15,17	1.00	2 (13%)
4	MAN	C	6	4	11,11,12	0.63	0	15,15,17	1.03	2 (13%)
3	NAG	F	1	3,1	14,14,15	0.25	0	17,19,21	0.62	0
3	NAG	F	2	3	14,14,15	0.35	0	17,19,21	0.40	0
3	NAG	H	1	3,1	14,14,15	0.18	0	17,19,21	0.46	0
3	NAG	H	2	3	14,14,15	0.33	0	17,19,21	0.37	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
4	NAG	C	1	4,2	-	0/6/23/26	0/1/1/1
4	NAG	C	2	4	-	2/6/23/26	0/1/1/1
4	BMA	C	3	4	-	2/2/19/22	0/1/1/1
4	MAN	C	4	4	-	1/2/19/22	0/1/1/1
4	MAN	C	5	4	-	1/2/19/22	0/1/1/1
4	MAN	C	6	4	-	1/2/19/22	0/1/1/1
3	NAG	F	1	3,1	-	2/6/23/26	0/1/1/1
3	NAG	F	2	3	-	2/6/23/26	0/1/1/1
3	NAG	H	1	3,1	-	0/6/23/26	0/1/1/1
3	NAG	H	2	3	-	2/6/23/26	0/1/1/1

There are no bond length outliers.

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	C	6	MAN	C1-O5-C5	2.57	115.63	112.19
4	C	5	MAN	C1-O5-C5	2.45	115.46	112.19
4	C	6	MAN	O2-C2-C3	-2.34	105.29	110.15
4	C	5	MAN	O2-C2-C3	-2.19	105.61	110.15
4	C	4	MAN	C1-O5-C5	2.13	115.05	112.19

There are no chirality outliers.

All (13) torsion outliers are listed below:

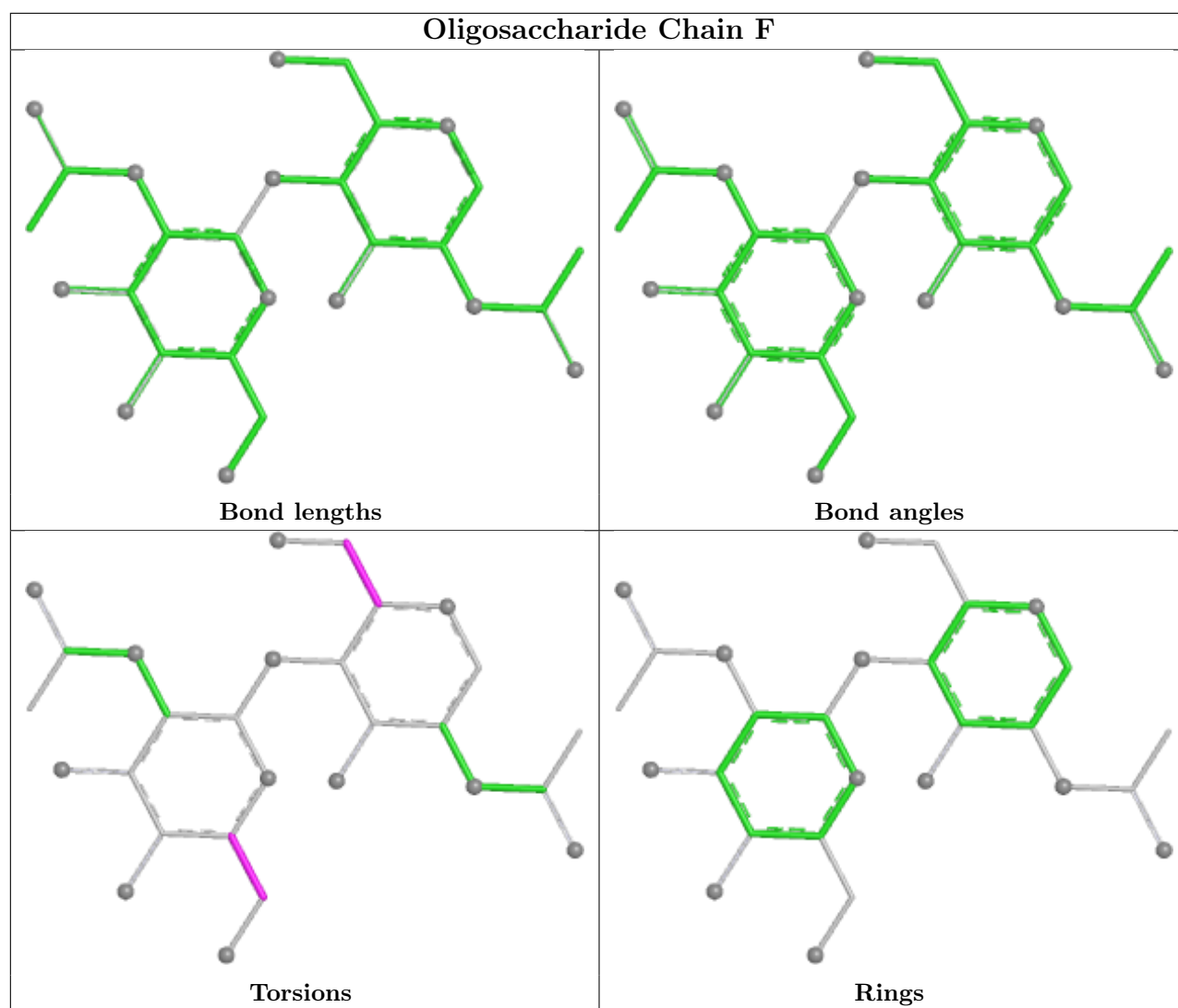
Mol	Chain	Res	Type	Atoms
4	C	3	BMA	O5-C5-C6-O6
3	H	2	NAG	O5-C5-C6-O6
3	H	2	NAG	C4-C5-C6-O6
4	C	3	BMA	C4-C5-C6-O6
3	F	2	NAG	C4-C5-C6-O6
4	C	2	NAG	C4-C5-C6-O6
3	F	1	NAG	C4-C5-C6-O6
3	F	2	NAG	O5-C5-C6-O6
3	F	1	NAG	O5-C5-C6-O6
4	C	2	NAG	O5-C5-C6-O6
4	C	6	MAN	C4-C5-C6-O6
4	C	4	MAN	C4-C5-C6-O6
4	C	5	MAN	C4-C5-C6-O6

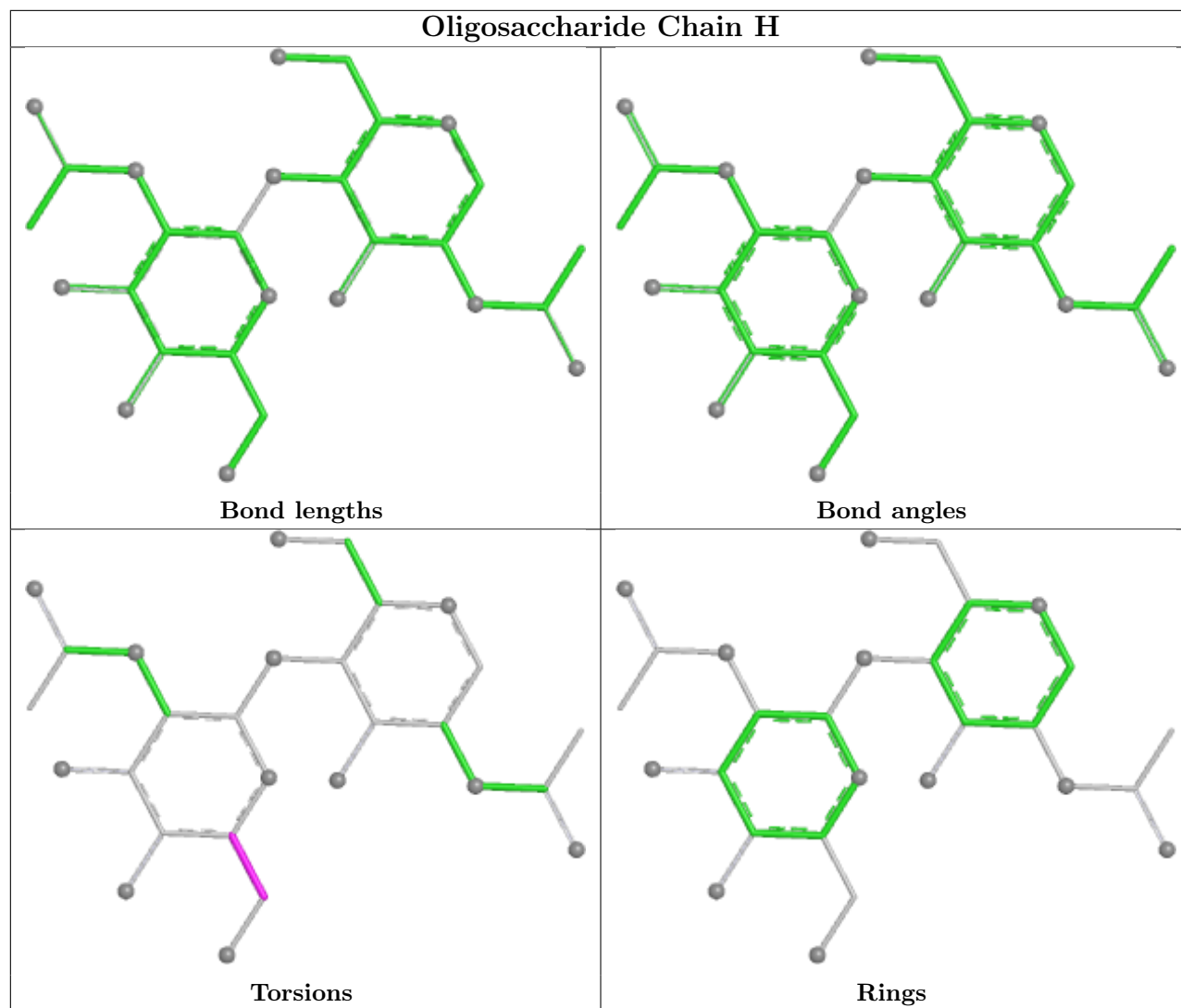
There are no ring outliers.

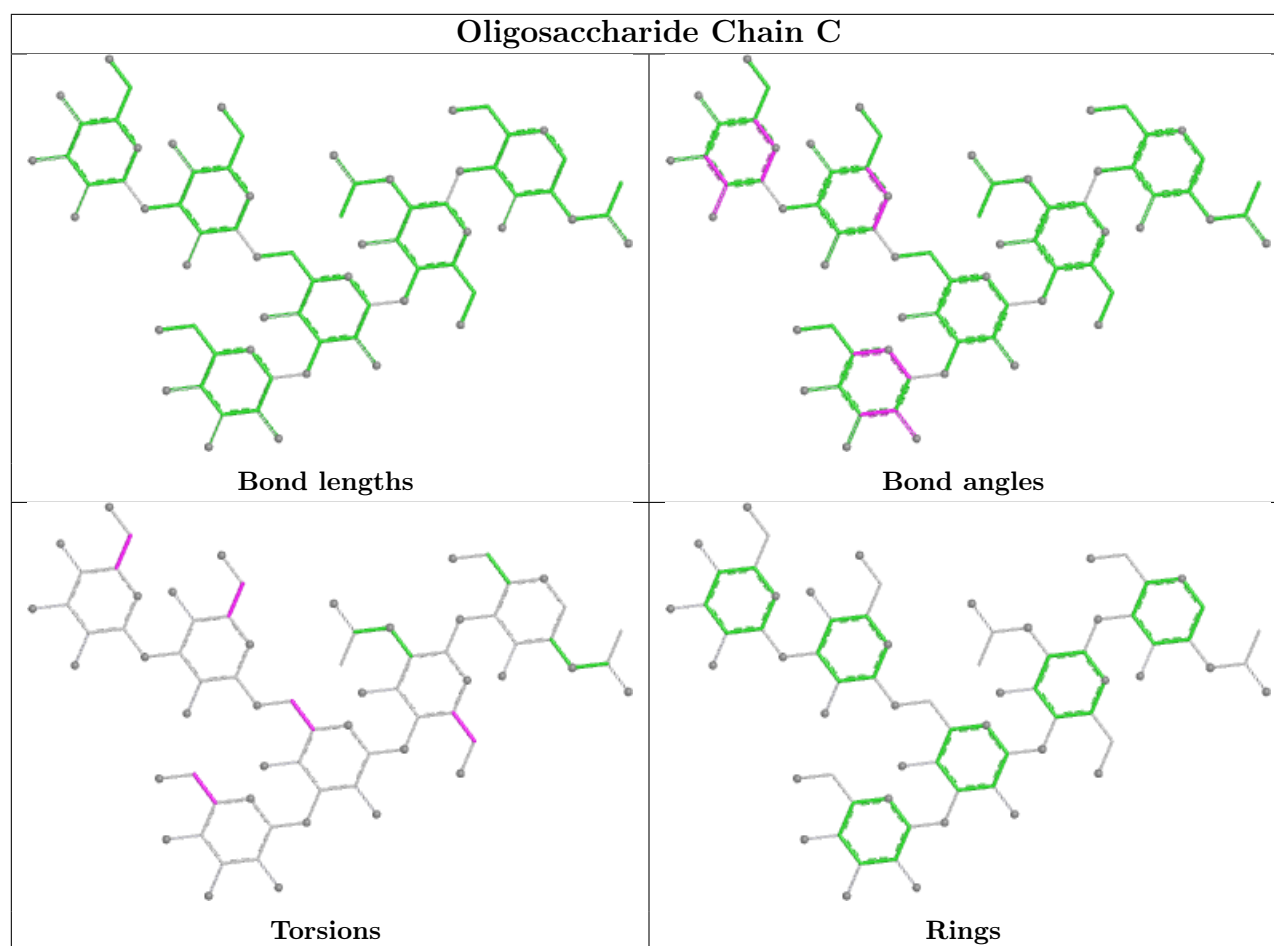
2 monomers are involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	F	2	NAG	1	0
3	F	1	NAG	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](#)

4 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	$\# Z > 2$
5	GOL	A	501	-	5,5,5	0.89	0	5,5,5	1.11	0
5	GOL	A	502	-	5,5,5	1.00	0	5,5,5	0.97	0
6	NAG	A	503	1	14,14,15	0.19	0	17,19,21	0.44	0
7	PO4	A	504	-	4,4,4	0.95	0	6,6,6	0.46	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
5	GOL	A	501	-	-	0/4/4/4	-
5	GOL	A	502	-	-	4/4/4/4	-
6	NAG	A	503	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 (5) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	A	502	GOL	O1-C1-C2-C3
5	A	502	GOL	C1-C2-C3-O3
5	A	502	GOL	O1-C1-C2-O2
5	A	502	GOL	O2-C2-C3-O3
6	A	503	NAG	C4-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](#)

The following chains have linkage breaks:

Mol	Chain	Number of breaks
1	A	1

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	A	293:THR	C	473:UNK	N	32.75

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	245/297 (82%)	0.45	27 (11%) 10 11	30, 49, 101, 137	1 (0%)
2	B	125/165 (75%)	0.26	9 (7%) 21 22	33, 45, 104, 147	0
All	All	370/462 (80%)	0.39	36 (9%) 13 14	30, 47, 104, 147	1 (0%)

All (36) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	626	ILE	7.4
1	A	198	PRO	6.2
1	A	288	TRP	5.4
1	A	283	THR	5.1
1	A	214	TYR	5.0
1	A	293	THR	4.9
2	B	525	ALA	4.4
1	A	46	SER	4.3
1	A	269	THR	4.3
1	A	279	PRO	4.2
2	B	524	GLY	4.1
2	B	623	ILE	4.1
1	A	281	ILE	3.9
1	A	278	ASN	3.9
2	B	527	ILE	3.5
2	B	526	ALA	3.3
1	A	194	PHE	3.3
1	A	50	LYS	3.2
1	A	191	LYS	3.2
1	A	282	ASP	3.2
1	A	197	HIS	3.2
1	A	47	ASP	3.0
1	A	51	LEU	2.7
1	A	280	GLU	2.7

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Mol	Chain	Res	Type	RSRZ
1	A	49	ASP	2.6
1	A	228	ASN	2.5
1	A	275	TRP	2.5
2	B	620	THR	2.4
1	A	48	VAL	2.4
1	A	277	VAL	2.3
2	B	621	ASP	2.2
1	A	229	GLU	2.2
1	A	270	THR	2.2
1	A	188	GLN	2.2
2	B	619	ILE	2.1
1	A	184	LEU	2.1

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

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

6.3 Carbohydrates ⓘ

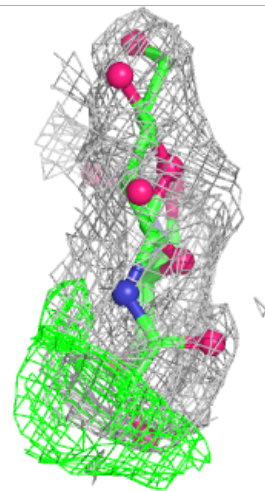
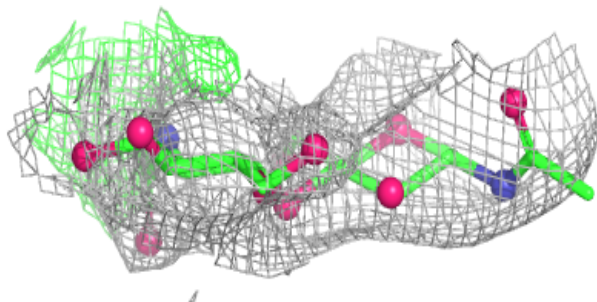
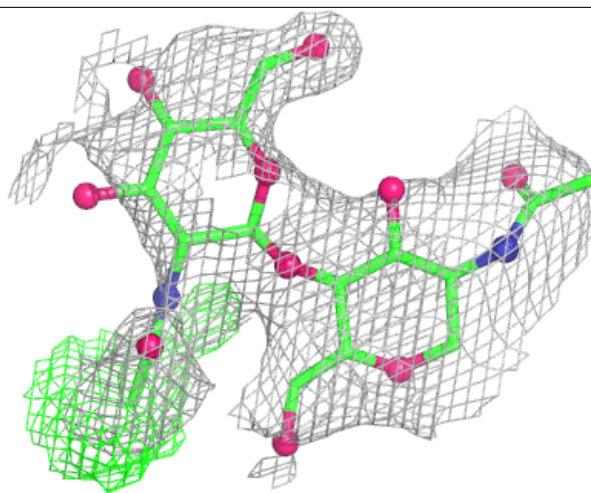
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists 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	NAG	F	1	14/15	-	-	91,105,119,120	0
3	NAG	F	2	14/15	-	-	99,136,139,142	0
3	NAG	H	1	14/15	-	-	61,75,85,95	0
3	NAG	H	2	14/15	-	-	85,108,116,119	0
4	NAG	C	2	14/15	0.23	0.23	44,67,80,86	0
4	NAG	C	1	14/15	0.86	0.13	35,43,52,54	0
4	BMA	C	3	11/12	-	-	71,84,95,101	0
4	MAN	C	4	11/12	-	-	91,99,106,114	0
4	MAN	C	5	11/12	-	-	90,112,123,128	0
4	MAN	C	6	11/12	-	-	86,99,115,117	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.

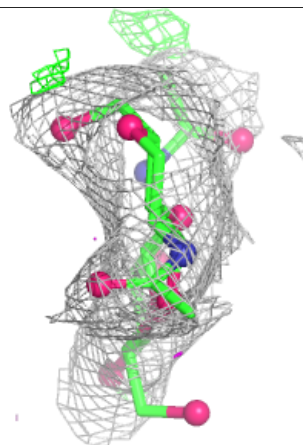
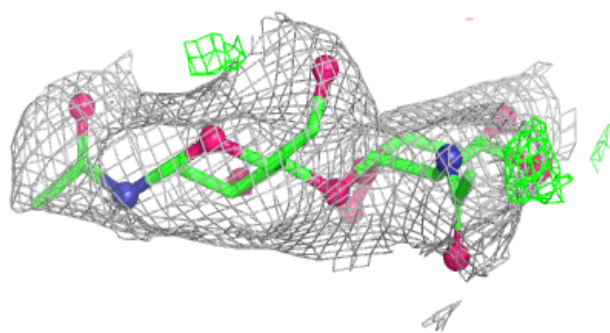
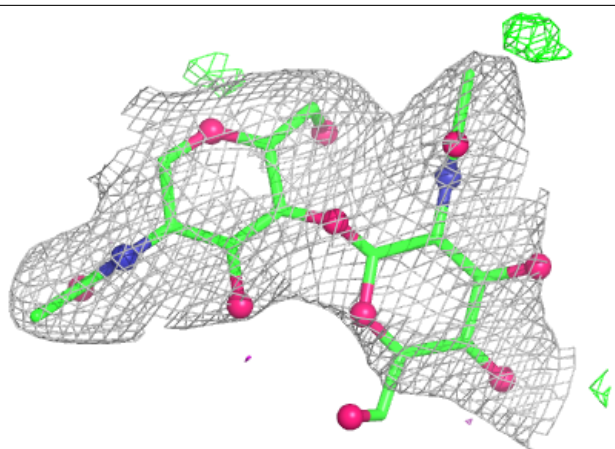
Electron density around Chain F:

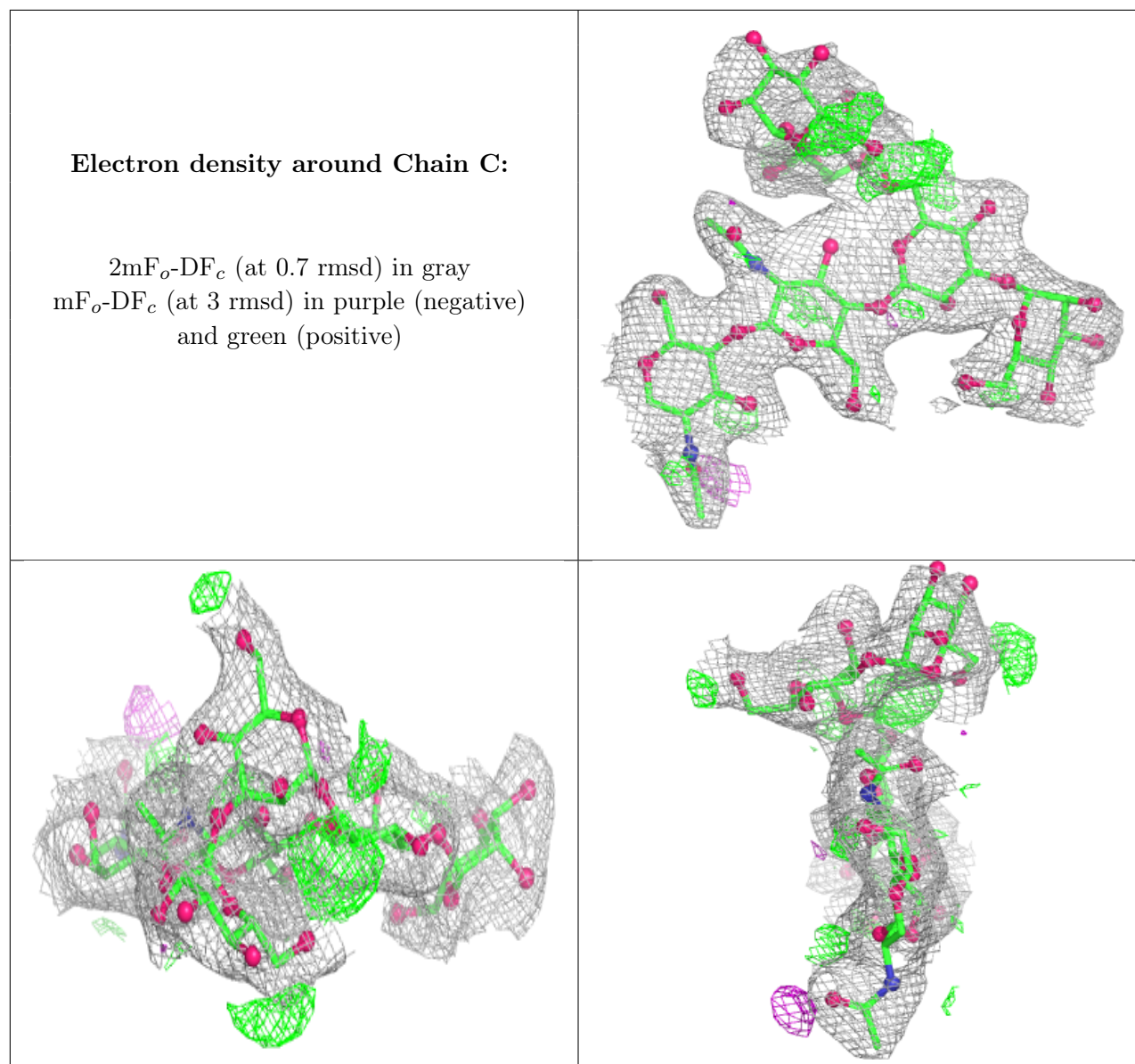
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around Chain H:

$2mF_o - DF_c$ (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(Å ²)	Q<0.9
6	NAG	A	503	14/15	0.53	0.18	85,133,144,147	0
5	GOL	A	502	6/6	0.71	0.23	73,82,104,107	0
7	PO4	A	504	5/5	0.82	0.13	72,87,112,129	0
5	GOL	A	501	6/6	0.83	0.20	64,72,88,92	0

6.5 Other polymers ⓘ

There are no such residues in this entry.