



# Full wwPDB X-ray Structure Validation Report ⓘ

Mar 10, 2025 – 01:06 PM EDT

PDB ID : 9MEV  
Title : Structure of H1H3:FluA20 Chimeric Antigen Complex  
Authors : Seraj, N.  
Deposited on : 2024-12-08  
Resolution : 2.06 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](mailto: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>.

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The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467  
Mogul : 2022.3.0, CSD as543be (2022)  
Xtriage (Phenix) : 1.21  
EDS : 3.0  
Percentile statistics : 20231227.v01 (using entries in the PDB archive December 27th 2023)  
CCP4 : 9.0.004 (Gargrove)  
Density-Fitness : 1.0.11  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.41.4

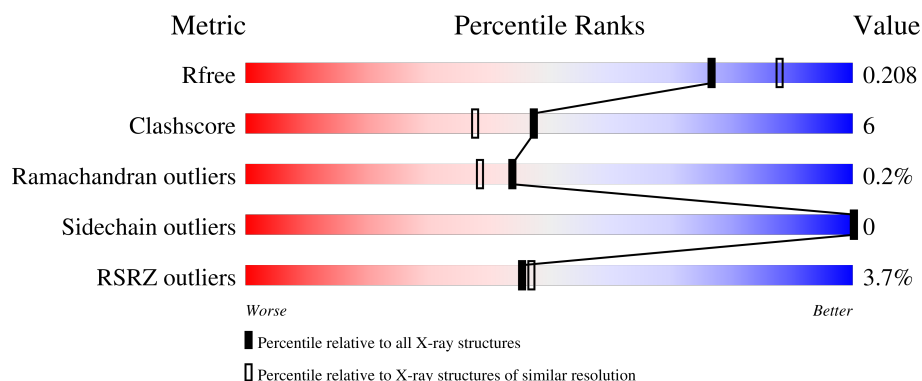
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.06 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))
$R_{free}$	164625	3436 (2.08-2.04)
Clashscore	180529	3661 (2.08-2.04)
Ramachandran outliers	177936	3649 (2.08-2.04)
Sidechain outliers	177891	3649 (2.08-2.04)
RSRZ outliers	164620	3436 (2.08-2.04)

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  $\geq 3$ , 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	B	227	<div> <div>5%</div> <div> <div></div> <div>89%</div> <div>5%</div> <div>7%</div> </div> </div>
2	H	235	<div> <div>4%</div> <div> <div></div> <div>77%</div> <div>19%</div> <div>.</div> </div> </div>
3	L	214	<div> <div>%</div> <div> <div></div> <div>90%</div> <div>9%</div> </div> </div>
4	A	2	<div> <div></div> <div>100%</div> </div>

## 2 Entry composition [i](#)

There are 7 unique types of molecules in this entry. The entry contains 5727 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 H1H3 HA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	B	212	Total	C	N	O	S	0	1	0
			1706	1086	290	326	4			

- Molecule 2 is a protein called FluA20 Heavy Chain Fab.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	H	227	Total	C	N	O	S	0	4	0
			1713	1080	279	347	7			

- Molecule 3 is a protein called FluA20 Light Chain Fab.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	L	213	Total	C	N	O	S	0	2	0
			1654	1038	281	330	5			

- Molecule 4 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
4	A	2	Total	C	N	O	0	0	0
			28	16	2	10			

- Molecule 5 is 2-acetamido-2-deoxy-beta-D-glucopyranose (three-letter code: NAG) (formula: C<sub>8</sub>H<sub>15</sub>NO<sub>6</sub>).



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

- Molecule 6 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula:  $C_2H_6O_2$ ).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	H	1	Total	C	O	0	0
			3	2	1		
6	L	1	Total	C	O	0	0
			4	2	2		
6	L	1	Total	C	O	0	0
			4	2	2		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	L	1	Total	C	O	0	0
			4	2	2		
6	L	1	Total	C	O	0	0
			4	2	2		
6	L	1	Total	C	O	0	0
			4	2	2		

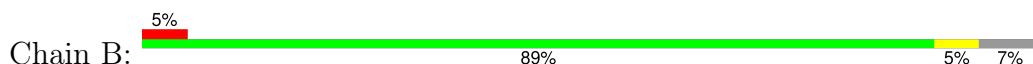
- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	B	165	Total	O	0	0
			165	165		
7	H	213	Total	O	0	0
			213	213		
7	L	211	Total	O	0	0
			211	211		

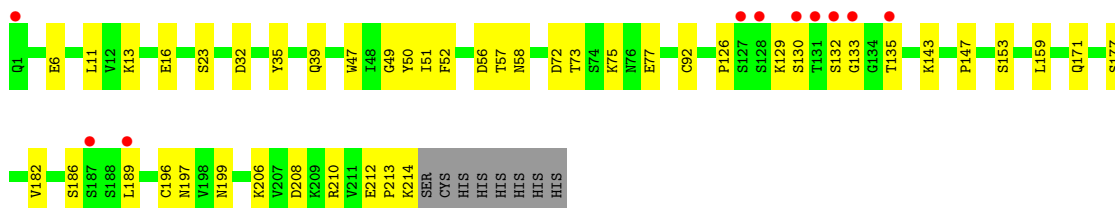
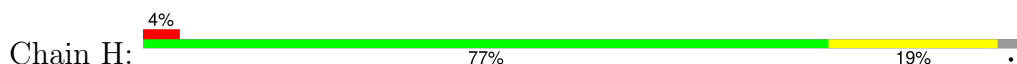
### 3 Residue-property plots [i](#)

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.

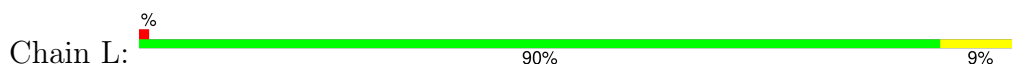
- Molecule 1: H1H3 HA



- Molecule 2: FluA20 Heavy Chain Fab



- Molecule 3: FluA20 Light Chain Fab



- Molecule 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, $\alpha$ , $\beta$ , $\gamma$	50.64Å 93.47Å 86.43Å 90.00° 93.48° 90.00°	Depositor
Resolution (Å)	44.82 – 2.06 44.82 – 2.06	Depositor EDS
% Data completeness (in resolution range)	72.9 (44.82-2.06) 92.0 (44.82-2.06)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.70 (at 2.05Å)	Xtriage
Refinement program	PHENIX 1.21.2_5419	Depositor
R, $R_{free}$	0.181 , 0.209 0.181 , 0.208	Depositor DCC
$R_{free}$ test set	47767 reflections (4.35%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	26.6	Xtriage
Anisotropy	0.049	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.35 , 53.9	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	5727	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	32.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.20% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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: EDO, 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	B	0.56	0/1760	0.59	0/2394
2	H	0.66	0/1768	0.70	0/2417
3	L	0.42	0/1697	0.56	0/2304
All	All	0.56	0/5225	0.62	0/7115

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	B	1706	0	1625	7	0
2	H	1713	0	1674	34	1
3	L	1654	0	1624	20	0
4	A	28	0	25	0	0
5	B	14	0	13	0	0
6	H	3	0	3	0	0
6	L	20	0	30	1	0
7	B	165	0	0	3	0
7	H	213	0	0	4	0
7	L	211	0	0	2	0
All	All	5727	0	4994	56	1



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 6.

All (56) 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:69:PRO:HA	1:B:149:LYS:HE2	1.60	0.82
3:L:33[B]:LEU:HD23	3:L:71:PHE:CD1	2.17	0.80
3:L:33[B]:LEU:HD11	3:L:88:CYS:HB2	1.66	0.78
3:L:80:PRO:HG2	6:L:304:EDO:H22	1.68	0.75
1:B:72:GLU:HB3	1:B:149:LYS:HE3	1.68	0.74
1:B:226:GLN:HG3	7:B:429:HOH:O	1.96	0.66
3:L:33[B]:LEU:HD12	3:L:34:ASN:N	2.10	0.65
3:L:169:LYS:HD2	7:L:496:HOH:O	1.96	0.64
3:L:33[B]:LEU:HD23	3:L:71:PHE:CG	2.33	0.64
2:H:132:SER:HB2	2:H:135:THR:OG1	2.01	0.60
2:H:126:PRO:HG2	2:H:213:PRO:HA	1.85	0.58
2:H:35:TYR:HB3	2:H:52:PHE:CZ	2.40	0.56
2:H:210:ARG:HH11	2:H:212:GLU:HG2	1.71	0.56
2:H:13:LYS:HB2	2:H:16:GLU:HG3	1.88	0.55
2:H:213:PRO:O	2:H:214:LYS:HB2	2.07	0.54
2:H:47:TRP:CH2	2:H:49:GLY:HA2	2.43	0.54
3:L:185:ASP:O	3:L:188:LYS:HG2	2.07	0.53
2:H:210:ARG:HD3	2:H:212:GLU:OE2	2.10	0.52
2:H:47:TRP:CE3	3:L:96:PRO:HD2	2.45	0.51
3:L:33[B]:LEU:HD12	3:L:34:ASN:H	1.74	0.51
2:H:47:TRP:CD2	3:L:96:PRO:HD2	2.47	0.50
2:H:186:SER:HA	2:H:189:LEU:HG	1.94	0.50
2:H:132:SER:HB3	7:H:433:HOH:O	2.11	0.50
3:L:33[A]:LEU:HD13	3:L:71:PHE:CD1	2.47	0.49
2:H:143:LYS:HG3	2:H:177[B]:SER:OG	2.13	0.49
2:H:210:ARG:HH11	2:H:212:GLU:CG	2.26	0.48
2:H:129:LYS:HB3	3:L:117:ILE:CG2	2.44	0.48
2:H:23:SER:HA	2:H:77[A]:GLU:HG2	1.95	0.48
2:H:72:ASP:HB3	2:H:77[B]:GLU:HG2	1.95	0.47
2:H:210:ARG:NH1	2:H:212:GLU:HG2	2.29	0.47
3:L:163:VAL:HG22	3:L:175:LEU:HD12	1.96	0.47
2:H:130:SER:HA	3:L:116:PHE:HD2	1.79	0.47
2:H:153:SER:OG	2:H:197:ASN:HB2	2.15	0.47
3:L:12[A]:SER:OG	3:L:107:LYS:HG2	2.15	0.47
2:H:47:TRP:CZ2	2:H:49:GLY:HA2	2.51	0.46
2:H:6:GLU:HG3	2:H:92:CYS:SG	2.57	0.45
2:H:143:LYS:NZ	2:H:171:GLN:OE1	2.50	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:L:207:LYS:HD2	7:L:415:HOH:O	2.16	0.45
2:H:73:THR:HG23	7:H:453:HOH:O	2.16	0.44
3:L:120:PRO:HD3	3:L:132:VAL:HG22	1.99	0.44
1:B:175:GLU:HG2	1:B:238:GLU:HA	2.00	0.44
2:H:51:ILE:HG13	2:H:57:THR:HG22	1.99	0.44
2:H:50:TYR:CE2	2:H:58:ASN:HB3	2.53	0.43
1:B:197:LYS:HD2	1:B:248:ASN:O	2.19	0.43
2:H:11:LEU:HG	2:H:147:PRO:HG3	1.99	0.43
2:H:72:ASP:OD2	2:H:75:LYS:HD2	2.19	0.42
3:L:29:ILE:HD11	3:L:33[B]:LEU:HD22	2.01	0.42
1:B:156:LYS:NZ	7:B:410:HOH:O	2.51	0.42
2:H:199:ASN:HB2	2:H:206:LYS:HE3	2.01	0.42
1:B:215:PRO:HG2	7:B:418:HOH:O	2.19	0.42
2:H:159:LEU:HD21	2:H:182:VAL:HG21	2.02	0.41
2:H:39:GLN:NE2	7:H:404:HOH:O	2.44	0.41
2:H:196:CYS:O	2:H:208:ASP:HA	2.20	0.41
2:H:129:LYS:HB3	3:L:117:ILE:HG22	2.02	0.40
2:H:32[A]:ASP:OD2	7:H:401:HOH:O	2.21	0.40
3:L:95:PRO:HA	3:L:96:PRO:HD3	2.01	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:56:ASP:OD2	2:H:133:GLY:O[2_645]	1.96	0.24

## 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	B	211/227 (93%)	203 (96%)	7 (3%)	1 (0%)	25 17

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
2	H	229/235 (97%)	224 (98%)	5 (2%)	0	100	100
3	L	213/214 (100%)	206 (97%)	7 (3%)	0	100	100
All	All	653/676 (97%)	633 (97%)	19 (3%)	1 (0%)	44	38

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	82	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	B	189/200 (94%)	189 (100%)	0	100	100
2	H	201/205 (98%)	201 (100%)	0	100	100
3	L	191/190 (100%)	191 (100%)	0	100	100
All	All	581/595 (98%)	581 (100%)	0	100	100

There are no protein residues with a non-rotameric sidechain to report.

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

2 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	A	1	1,4	14,14,15	0.41	0	17,19,21	0.66	0
4	NAG	A	2	4	14,14,15	0.40	0	17,19,21	0.95	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	A	1	1,4	-	0/6/23/26	0/1/1/1
4	NAG	A	2	4	-	0/6/23/26	0/1/1/1

There are no bond length outliers.

There are no bond angle outliers.

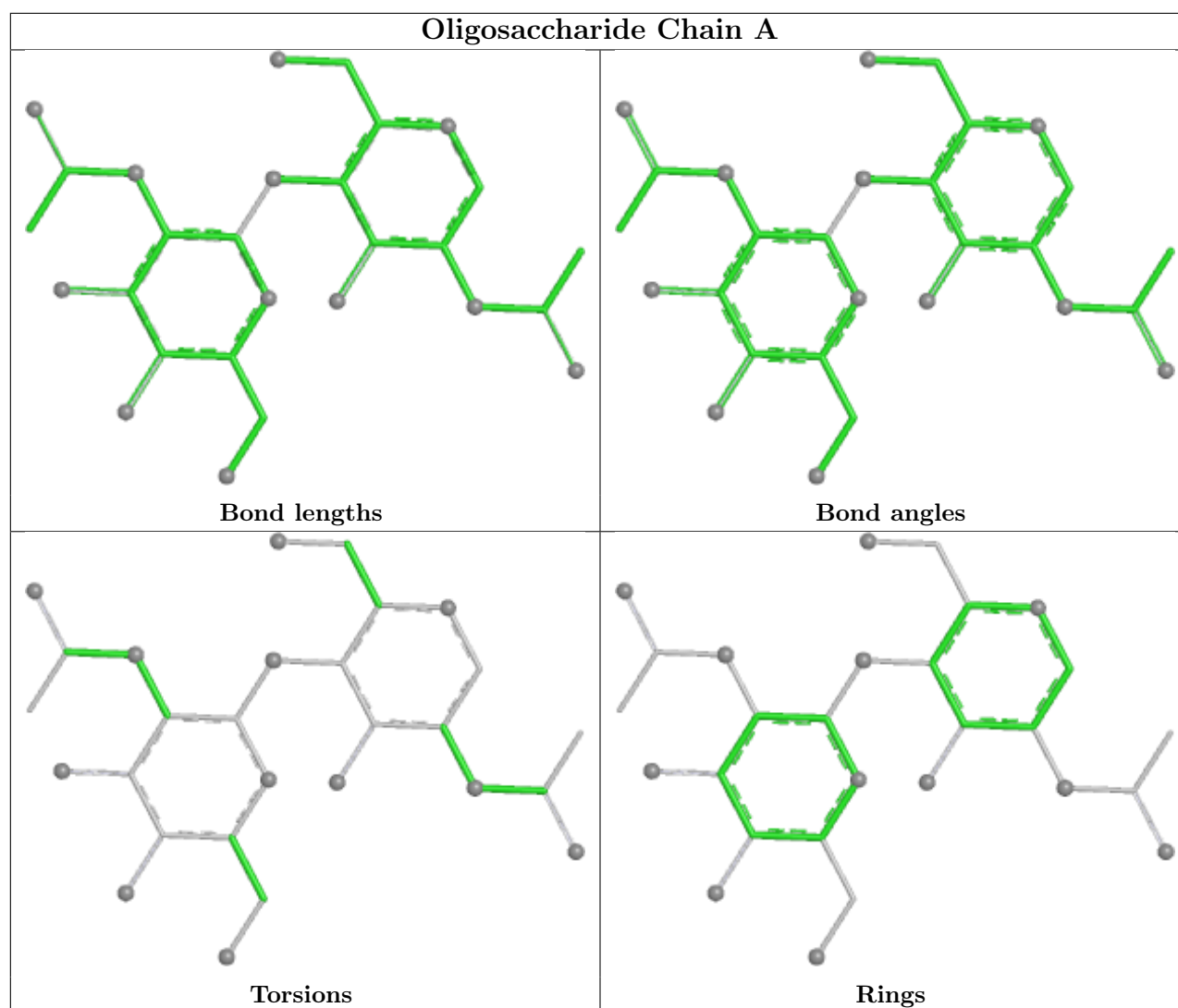
There are no chirality outliers.

There are no torsion outliers.

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](#)

7 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	$\# Z  > 2$
6	EDO	L	301	-	3,3,3	0.38	0	2,2,2	0.24	0
6	EDO	L	303	-	3,3,3	0.39	0	2,2,2	0.27	0
5	NAG	B	301	1	14,14,15	0.36	0	17,19,21	0.72	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
6	EDO	L	305	-	3,3,3	0.35	0	2,2,2	0.37	0
6	EDO	L	302	-	3,3,3	0.42	0	2,2,2	0.17	0
6	EDO	H	301	-	2,2,3	0.42	0	1,1,2	0.33	0
6	EDO	L	304	-	3,3,3	0.33	0	2,2,2	0.50	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
6	EDO	L	301	-	-	0/1/1/1	-
6	EDO	L	303	-	-	1/1/1/1	-
5	NAG	B	301	1	-	0/6/23/26	0/1/1/1
6	EDO	L	305	-	-	1/1/1/1	-
6	EDO	L	302	-	-	0/1/1/1	-
6	EDO	L	304	-	-	0/1/1/1	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
6	L	305	EDO	O1-C1-C2-O2
6	L	303	EDO	O1-C1-C2-O2

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	L	304	EDO	1	0

## 5.7 Other polymers ⓘ

There are no such residues in this entry.

## 5.8 Polymer linkage issues ⓘ

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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	B	212/227 (93%)	0.20	11 (5%) 34 35	17, 33, 55, 76	1 (0%)
2	H	227/235 (96%)	-0.01	10 (4%) 39 42	13, 26, 56, 75	4 (1%)
3	L	213/214 (99%)	-0.13	3 (1%) 73 75	12, 25, 54, 74	2 (0%)
All	All	652/676 (96%)	0.02	24 (3%) 45 47	12, 28, 56, 76	7 (1%)

All (24) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	H	132	SER	4.7
1	B	75	ILE	4.2
2	H	130	SER	3.9
2	H	131	THR	3.5
2	H	133	GLY	2.9
2	H	128	SER	2.8
1	B	134	GLY	2.8
1	B	240	GLY	2.8
3	L	76	ARG	2.8
2	H	127	SER	2.8
2	H	187	SER	2.7
1	B	133	ASN	2.6
1	B	135	GLY	2.6
2	H	135	THR	2.6
3	L	1	ASP	2.5
1	B	74	LEU	2.5
2	H	189	LEU	2.5
3	L	152	ASN	2.5
1	B	169	ALA	2.4
1	B	263	GLY	2.4
1	B	132	GLN	2.2
1	B	73	LEU	2.2
1	B	171	ASN	2.1

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Mol	Chain	Res	Type	RSRZ
2	H	1	GLN	2.1

## 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](#)

SUGAR-RSR INFOmissingINFO

## 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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
5	NAG	B	301	14/15	0.66	0.16	51,67,81,88	0
6	EDO	H	301	3/4	0.68	0.22	26,26,27,35	0
6	EDO	L	302	4/4	0.81	0.17	43,49,51,53	0
6	EDO	L	304	4/4	0.86	0.14	24,26,27,28	0
6	EDO	L	301	4/4	0.92	0.12	27,31,36,37	0
6	EDO	L	305	4/4	0.92	0.11	36,41,43,44	0
6	EDO	L	303	4/4	0.97	0.06	16,23,26,30	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.