



Full wwPDB X-ray Structure Validation Report ⓘ

Aug 21, 2020 – 11:02 PM BST

PDB ID : 5N0B
Title : Crystal structure of the tetanus neurotoxin in complex with GD1a
Authors : Masuyer, G.; Conrad, J.; Stenmark, P.
Deposited on : 2017-02-02
Resolution : 2.60 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 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 : 2.13.1
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) : 2.13.1

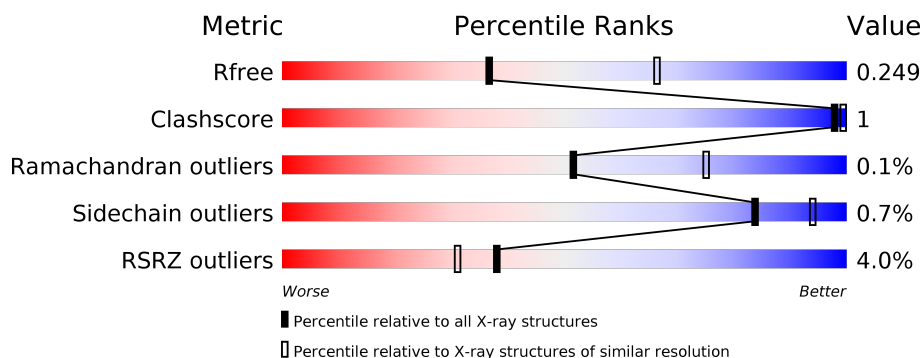
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.60 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	3163 (2.60-2.60)
Clashscore	141614	3518 (2.60-2.60)
Ramachandran outliers	138981	3455 (2.60-2.60)
Sidechain outliers	138945	3455 (2.60-2.60)
RSRZ outliers	127900	3104 (2.60-2.60)

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 on 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	1335	<div> <div>4%</div> <div>93%</div> <div>.</div> <div>.</div> </div>
2	B	6	<div> <div>17%</div> <div>83%</div> </div>

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 10580 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.

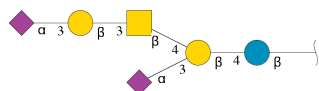
- Molecule 1 is a protein called Tetanus toxin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	1280	Total	C	N	O	S	0	0	0
			10348	6635	1689	1989	35			

There are 22 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-19	MET	-	initiating methionine	UNP P04958
A	-18	GLY	-	expression tag	UNP P04958
A	-17	SER	-	expression tag	UNP P04958
A	-16	SER	-	expression tag	UNP P04958
A	-15	HIS	-	expression tag	UNP P04958
A	-14	HIS	-	expression tag	UNP P04958
A	-13	HIS	-	expression tag	UNP P04958
A	-12	HIS	-	expression tag	UNP P04958
A	-11	HIS	-	expression tag	UNP P04958
A	-10	HIS	-	expression tag	UNP P04958
A	-9	SER	-	expression tag	UNP P04958
A	-8	SER	-	expression tag	UNP P04958
A	-7	GLY	-	expression tag	UNP P04958
A	-6	LEU	-	expression tag	UNP P04958
A	-5	VAL	-	expression tag	UNP P04958
A	-4	PRO	-	expression tag	UNP P04958
A	-3	ARG	-	expression tag	UNP P04958
A	-2	GLY	-	expression tag	UNP P04958
A	-1	SER	-	expression tag	UNP P04958
A	0	HIS	-	expression tag	UNP P04958
A	372	ALA	ARG	engineered mutation	UNP P04958
A	375	PHE	TYR	engineered mutation	UNP P04958

- Molecule 2 is an oligosaccharide called N-acetyl-alpha-neuraminic acid-(2-3)-beta-D-galactopyranose-(1-3)-2-acetamido-2-deoxy-beta-D-galactopyranose-(1-4)-[N-acetyl-alpha-neuraminic acid-(2-3)]beta-D-galactopyranose-(1-4)-beta-D-glucopyranose.

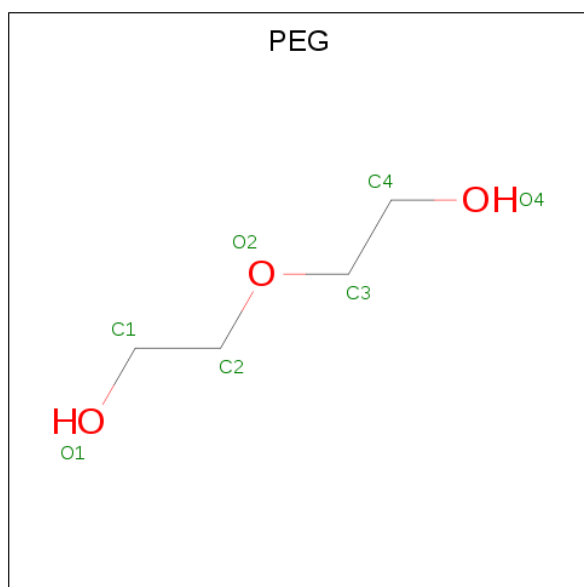


Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	B	6	Total	C	N	O	0	0	0
			88	48	3	37			

- Molecule 3 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	Zn	0	0
			1	1		

- Molecule 4 is DI(HYDROXYETHYL)ETHER (three-letter code: PEG) (formula: C₄H₁₀O₃).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			7	4	3		

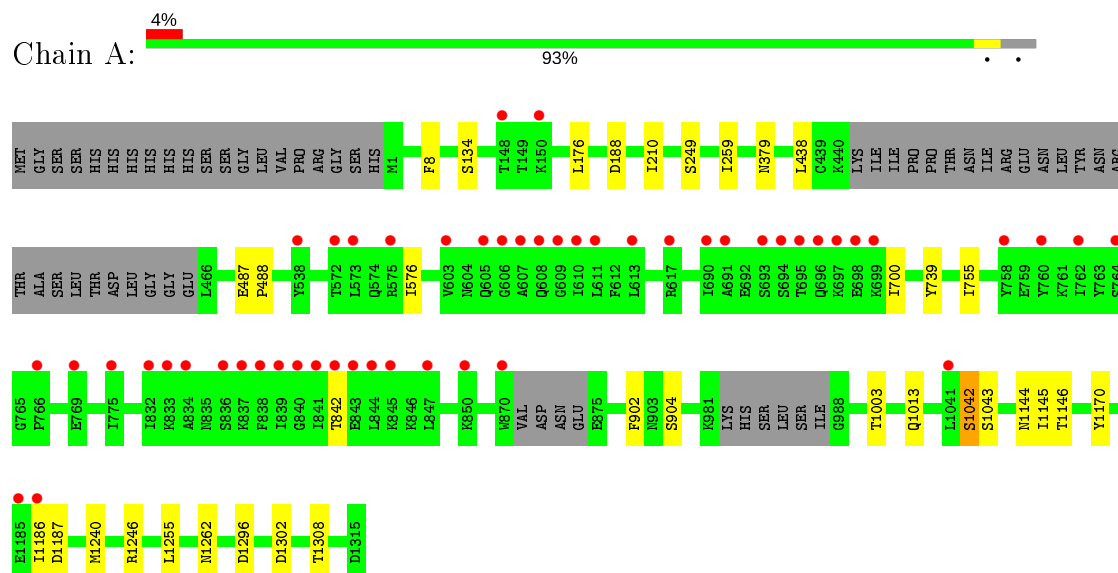
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	136	Total	O	0	0
			136	136		

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: Tetanus toxin



- Molecule 2: N-acetyl-alpha-neuraminic acid-(2-3)-beta-D-galactopyranose-(1-3)-2-acetamido-2-deoxy-beta-D-galactopyranose-(1-4)-[N-acetyl-alpha-neuraminic acid-(2-3)]beta-D-galactopyranose-(1-4)-beta-D-glucopyranose



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	152.69Å 136.84Å 92.63Å 90.00° 90.07° 90.00°	Depositor
Resolution (Å)	68.50 – 2.60 68.52 – 2.60	Depositor EDS
% Data completeness (in resolution range)	99.8 (68.50-2.60) 99.8 (68.52-2.60)	Depositor EDS
R_{merge}	0.13	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.74 (at 2.62Å)	Xtriage
Refinement program	REFMAC 5.8.0158	Depositor
R, R_{free}	0.203 , 0.247 0.208 , 0.249	Depositor DCC
R_{free} test set	2862 reflections (4.90%)	wwPDB-VP
Wilson B-factor (Å ²)	51.9	Xtriage
Anisotropy	0.197	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 50.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	0.025 for -h,-k,l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	10580	wwPDB-VP
Average B, all atoms (Å ²)	66.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.66% 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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: ZN, BGC, NGA, SIA, GAL, PEG

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	0.40	0/10568	0.58	0/14317

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 [i](#)

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	10348	0	10277	15	0
2	B	88	0	74	0	0
3	A	1	0	0	0	0
4	A	7	0	10	1	0
5	A	136	0	0	0	0
All	All	10580	0	10361	15	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 1.

All (15) 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:1145:ILE:O	1:A:1146:THR:HG22	2.01	0.59
1:A:1246:ARG:NH2	1:A:1296:ASP:O	2.38	0.56
1:A:1003:THR:HG23	1:A:1013:GLN:HE21	1.76	0.50
1:A:1240:MET:CE	1:A:1255:LEU:HB3	2.42	0.49
1:A:210:ILE:HD11	1:A:739:TYR:CE2	2.47	0.49
1:A:1170:TYR:CE1	1:A:1308:THR:HA	2.47	0.48
1:A:259:ILE:HG21	1:A:438:LEU:HD21	1.99	0.45
1:A:1246:ARG:NH1	1:A:1302:ASP:OD2	2.50	0.45
1:A:379:ASN:HD21	4:A:1402:PEG:H22	1.83	0.44
1:A:700:ILE:O	1:A:700:ILE:HG23	2.19	0.43
1:A:134:SER:HB3	1:A:176:LEU:HD21	2.00	0.42
1:A:576:ILE:HG22	1:A:755:ILE:HD11	2.01	0.42
1:A:487:GLU:N	1:A:488:PRO:CD	2.83	0.41
1:A:902:PHE:O	1:A:904:SER:N	2.53	0.41
1:A:1042:SER:OG	1:A:1043:SER:N	2.52	0.41

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	1272/1335 (95%)	1230 (97%)	41 (3%)	1 (0%)	51 75

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	1042	SER

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	1167/1216 (96%)	1159 (99%)	8 (1%)	84	94

All (8) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	8	PHE
1	A	188	ASP
1	A	249	SER
1	A	842	THR
1	A	1144	ASN
1	A	1186	ILE
1	A	1187	ASP
1	A	1262	ASN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (14) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	304	ASN
1	A	379	ASN
1	A	588	ASN
1	A	629	GLN
1	A	780	ASN
1	A	903	ASN
1	A	948	ASN
1	A	972	ASN
1	A	998	ASN
1	A	999	ASN
1	A	1013	GLN
1	A	1080	ASN
1	A	1272	ASN
1	A	1280	ASN

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 ⓘ

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	BGC	B	1	2	12,12,12	0.76	0	17,17,17	1.52	5 (29%)
2	GAL	B	2	2	11,11,12	0.41	0	15,15,17	1.56	2 (13%)
2	NGA	B	3	2	14,14,15	0.30	0	17,19,21	0.85	0
2	GAL	B	4	2	11,11,12	0.28	0	15,15,17	0.88	1 (6%)
2	SIA	B	5	2	17,20,21	0.61	0	21,28,31	1.06	1 (4%)
2	SIA	B	6	2	17,20,21	0.54	0	21,28,31	1.22	2 (9%)

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	BGC	B	1	2	-	2/2/22/22	0/1/1/1
2	GAL	B	2	2	-	2/2/19/22	0/1/1/1
2	NGA	B	3	2	-	2/6/23/26	0/1/1/1
2	GAL	B	4	2	-	2/2/19/22	0/1/1/1
2	SIA	B	5	2	-	4/14/34/38	0/1/1/1
2	SIA	B	6	2	-	2/14/34/38	0/1/1/1

There are no bond length outliers.

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	2	GAL	C1-O5-C5	4.34	118.07	112.19
2	B	6	SIA	C6-O6-C2	2.97	117.69	111.34
2	B	5	SIA	C6-O6-C2	2.86	117.45	111.34
2	B	1	BGC	C1-O5-C5	2.81	118.97	113.66
2	B	1	BGC	O5-C1-C2	2.80	115.28	110.28
2	B	6	SIA	C11-C10-N5	2.80	120.83	116.10
2	B	2	GAL	C1-C2-C3	2.68	112.96	109.67
2	B	1	BGC	C1-C2-C3	2.50	115.50	110.31
2	B	1	BGC	O4-C4-C3	2.25	115.54	110.35
2	B	4	GAL	O5-C5-C6	2.04	110.40	107.20
2	B	1	BGC	O4-C4-C5	2.01	114.29	109.30

There are no chirality outliers.

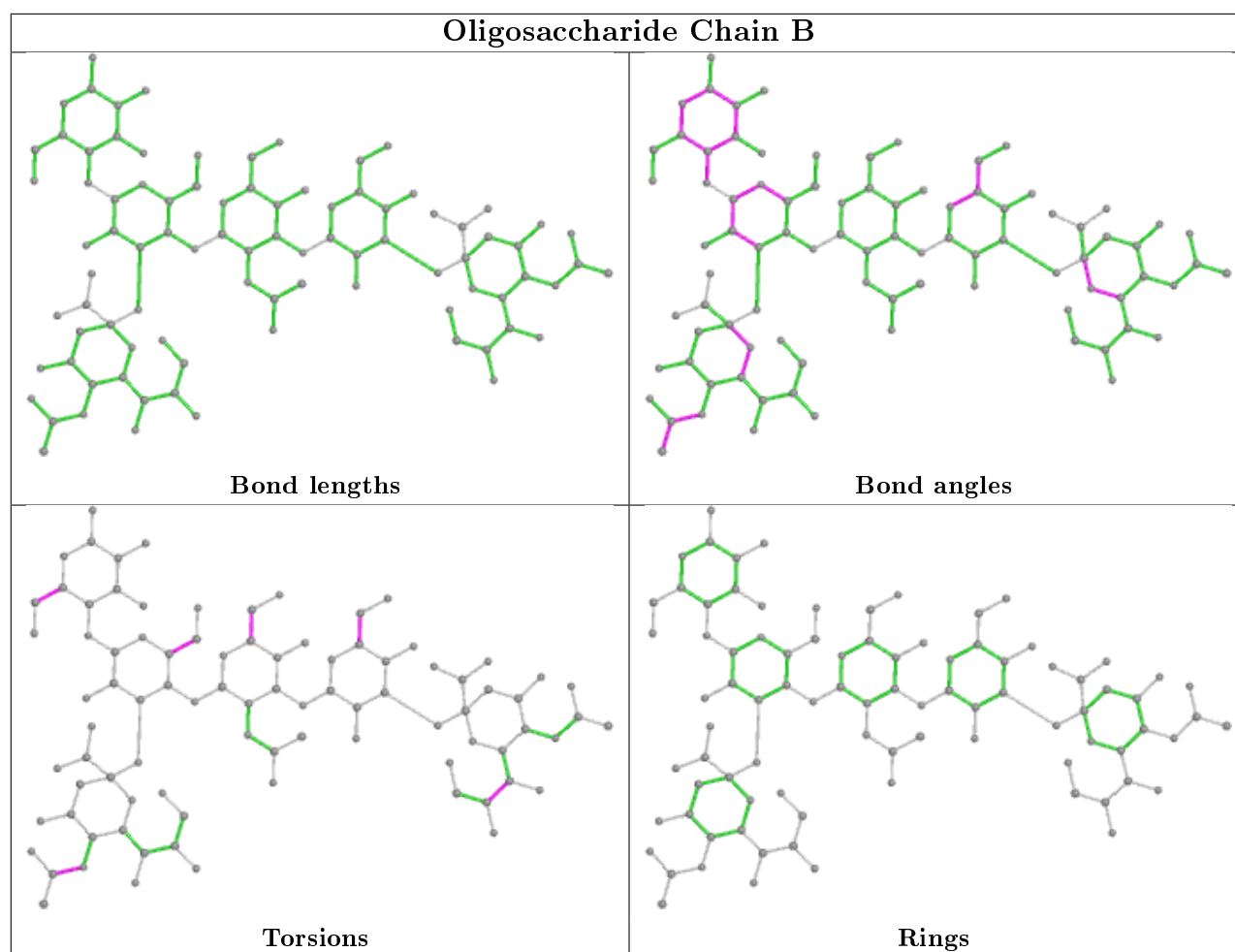
All (14) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	B	5	SIA	C6-C7-C8-C9
2	B	5	SIA	C6-C7-C8-O8
2	B	3	NGA	O5-C5-C6-O6
2	B	5	SIA	O7-C7-C8-O8
2	B	2	GAL	C4-C5-C6-O6
2	B	6	SIA	C11-C10-N5-C5
2	B	6	SIA	O10-C10-N5-C5
2	B	3	NGA	C4-C5-C6-O6
2	B	2	GAL	O5-C5-C6-O6
2	B	5	SIA	O7-C7-C8-C9
2	B	4	GAL	C4-C5-C6-O6
2	B	1	BGC	C4-C5-C6-O6
2	B	4	GAL	O5-C5-C6-O6
2	B	1	BGC	O5-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](#)

Of 2 ligands modelled in this entry, 1 is monoatomic - leaving 1 for Mogul analysis.

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	PEG	A	1402	-	6,6,6	0.47	0	5,5,5	0.33	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	PEG	A	1402	-	-	2/4/4/4	-

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
4	A	1402	PEG	O2-C3-C4-O4
4	A	1402	PEG	C1-C2-O2-C3

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	1402	PEG	1	0

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	1280/1335 (95%)	0.00	51 (3%) 38 31	31, 60, 118, 193	0

All (51) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	841	ILE	11.7
1	A	842	THR	10.3
1	A	698	GLU	8.5
1	A	839	ILE	8.4
1	A	843	GLU	7.5
1	A	696	GLN	6.4
1	A	608	GLN	6.4
1	A	613	LEU	6.1
1	A	609	GLY	5.2
1	A	575	ARG	5.1
1	A	840	GLY	4.5
1	A	573	LEU	4.5
1	A	693	SER	4.5
1	A	1185	GLU	4.1
1	A	847	LEU	3.9
1	A	697	LYS	3.8
1	A	837	LYS	3.8
1	A	538	TYR	3.6
1	A	695	THR	3.5
1	A	691	ALA	3.5
1	A	758	TYR	3.4
1	A	606	GLY	3.4
1	A	844	LEU	3.4
1	A	762	ILE	3.3
1	A	838	PHE	3.1
1	A	764	SER	3.1
1	A	572	THR	3.0

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Mol	Chain	Res	Type	RSRZ
1	A	766	PRO	3.0
1	A	1041	LEU	2.9
1	A	836	SER	2.9
1	A	690	ILE	2.8
1	A	832	ILE	2.7
1	A	834	ALA	2.6
1	A	760	TYR	2.6
1	A	603	VAL	2.4
1	A	694	SER	2.4
1	A	775	ILE	2.4
1	A	850	LYS	2.4
1	A	769	GLU	2.3
1	A	870	TRP	2.3
1	A	617	ARG	2.3
1	A	610	ILE	2.3
1	A	605	GLN	2.3
1	A	148	THR	2.2
1	A	833	LYS	2.2
1	A	611	LEU	2.2
1	A	150	LYS	2.1
1	A	845	LYS	2.1
1	A	1186	ILE	2.1
1	A	607	ALA	2.1
1	A	699	LYS	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.

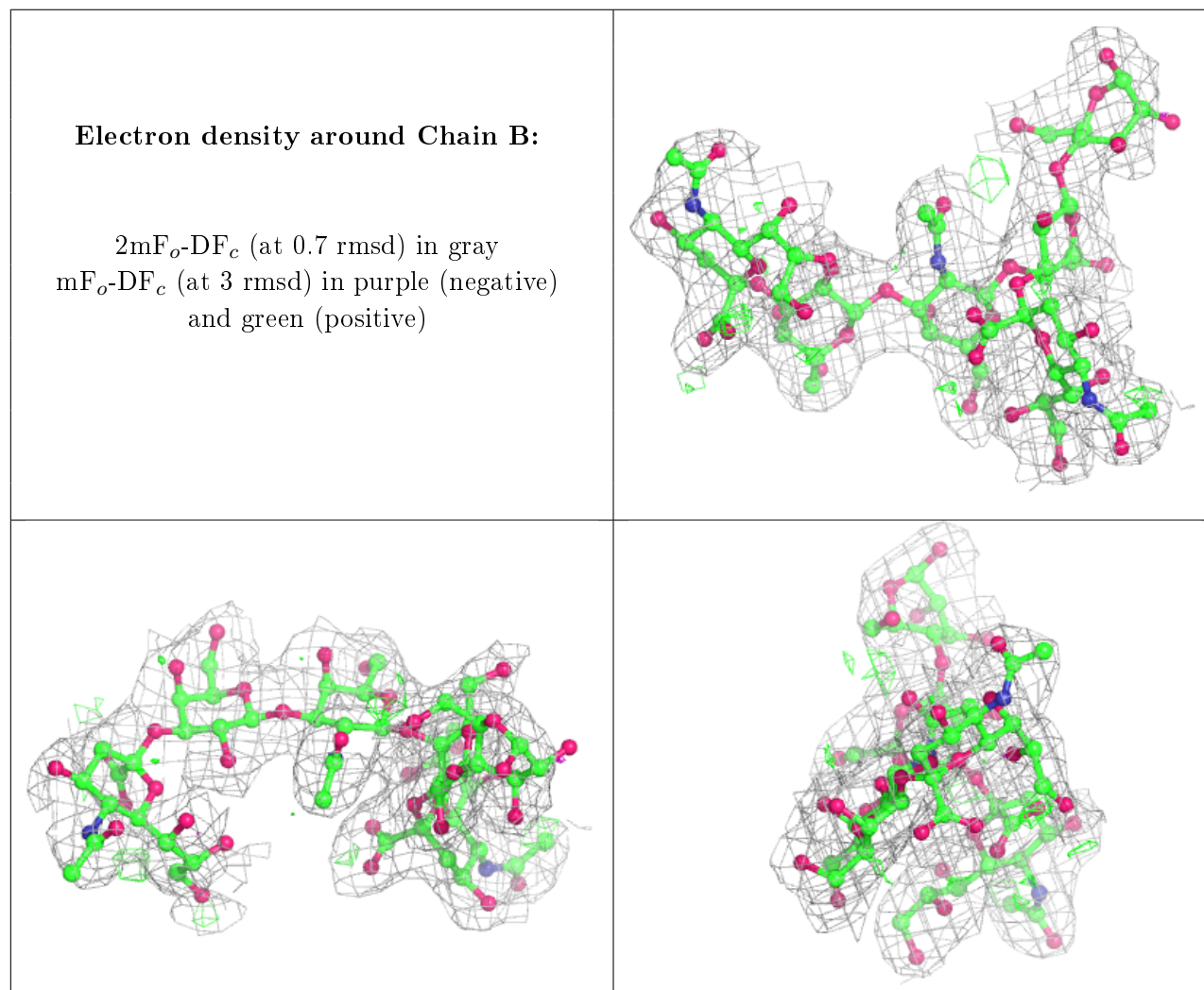
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	BGC	B	1	12/12	0.88	0.19	73,85,89,93	0
2	SIA	B	5	20/21	0.89	0.15	70,81,92,94	0
2	SIA	B	6	20/21	0.94	0.14	52,58,59,60	0
2	GAL	B	4	11/12	0.97	0.13	50,52,53,57	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
2	GAL	B	2	11/12	0.97	0.12	55,60,64,68	0
2	NGA	B	3	14/15	0.98	0.14	44,48,50,52	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(\AA^2)	Q<0.9
4	PEG	A	1402	7/7	0.71	0.23	81,86,88,88	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
3	ZN	A	1401	1/1	1.00	0.13	70,70,70,70	0

6.5 Other polymers [i](#)

There are no such residues in this entry.