



wwPDB X-ray Structure Validation Summary Report ⓘ

Apr 25, 2026 – 08:55 AM EDT

PDB ID : 5B0S / pdb_00005b0s
Title : Beta-1,2-Mannobiose phosphorylase from *Listeria innocua* - beta-1,2-mannotriose complex
Authors : Tsuda, T.; Arakawa, T.; Fushinobu, S.
Deposited on : 2015-11-02
Resolution : 2.10 Å(reported)

This is a wwPDB X-ray Structure Validation Summary 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

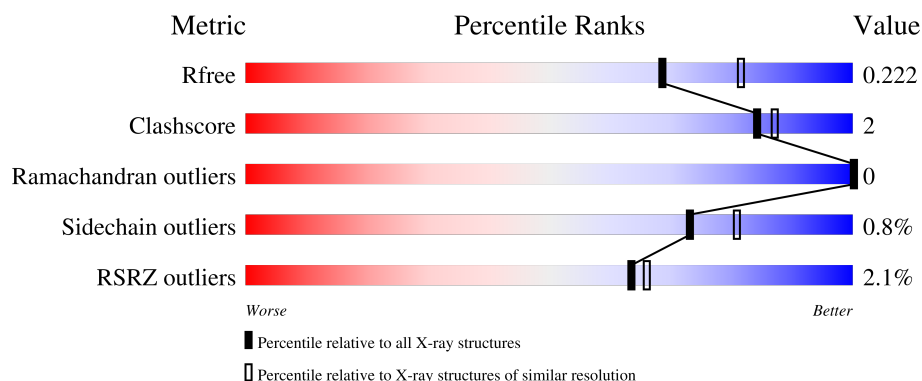
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.10 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	6658 (2.10-2.10)
Clashscore	190562	7164 (2.10-2.10)
Ramachandran outliers	187476	7099 (2.10-2.10)
Sidechain outliers	187428	7100 (2.10-2.10)
RSRZ outliers	180081	6662 (2.10-2.10)

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	363	2% 90% 7%
1	B	363	2% 90% 7%
2	C	3	67% 33%
2	D	3	33% 67%

2 Entry composition [i](#)

There are 6 unique types of molecules in this entry. The entry contains 6149 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.

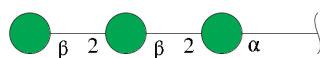
- Molecule 1 is a protein called Lin0857 protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	354	Total	C	N	O	S	0	0	0
			2815	1802	462	542	9			
1	B	355	Total	C	N	O	S	0	0	0
			2824	1808	464	543	9			

There are 16 discrepancies between the modelled and reference sequences:

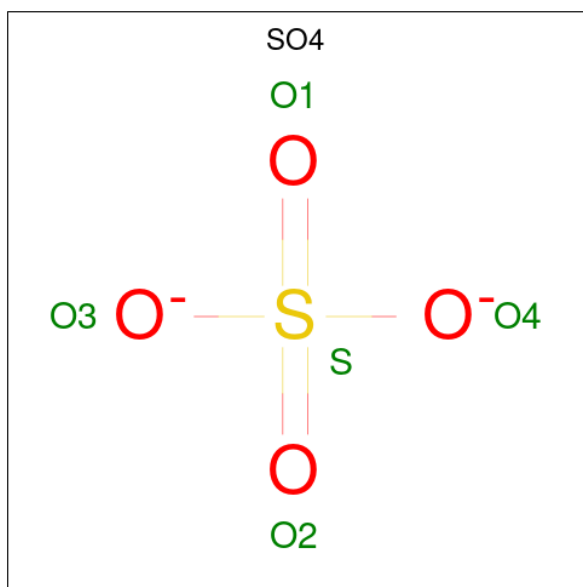
Chain	Residue	Modelled	Actual	Comment	Reference
A	356	LEU	-	expression tag	UNP Q92DF6
A	357	GLU	-	expression tag	UNP Q92DF6
A	358	HIS	-	expression tag	UNP Q92DF6
A	359	HIS	-	expression tag	UNP Q92DF6
A	360	HIS	-	expression tag	UNP Q92DF6
A	361	HIS	-	expression tag	UNP Q92DF6
A	362	HIS	-	expression tag	UNP Q92DF6
A	363	HIS	-	expression tag	UNP Q92DF6
B	356	LEU	-	expression tag	UNP Q92DF6
B	357	GLU	-	expression tag	UNP Q92DF6
B	358	HIS	-	expression tag	UNP Q92DF6
B	359	HIS	-	expression tag	UNP Q92DF6
B	360	HIS	-	expression tag	UNP Q92DF6
B	361	HIS	-	expression tag	UNP Q92DF6
B	362	HIS	-	expression tag	UNP Q92DF6
B	363	HIS	-	expression tag	UNP Q92DF6

- Molecule 2 is an oligosaccharide called beta-D-mannopyranose-(1-2)-beta-D-mannopyranose-(1-2)-alpha-D-mannopyranose.



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace
2	C	3	Total	C	O	0	0	0
			34	18	16			
2	D	3	Total	C	O	0	0	0
			34	18	16			

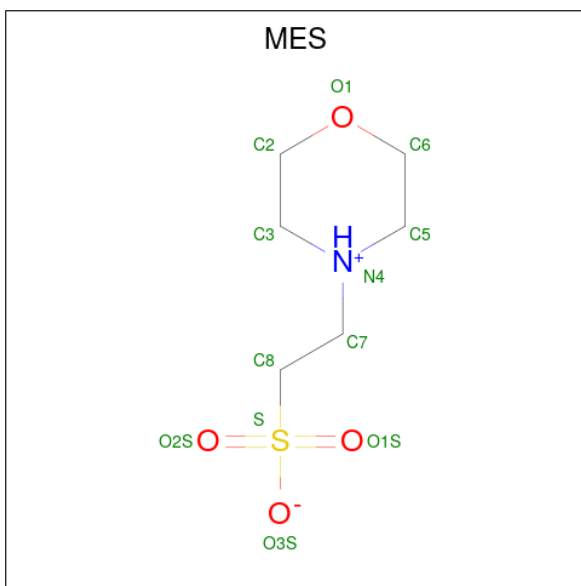
- Molecule 3 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	O	S	0	0
			5	4	1		
3	A	1	Total	O	S	0	0
			5	4	1		
3	A	1	Total	O	S	0	0
			5	4	1		
3	A	1	Total	O	S	0	0
			5	4	1		
3	A	1	Total	O	S	0	0
			5	4	1		
3	B	1	Total	O	S	0	0
			5	4	1		
3	B	1	Total	O	S	0	0
			5	4	1		
3	B	1	Total	O	S	0	0
			5	4	1		
3	B	1	Total	O	S	0	0
			5	4	1		

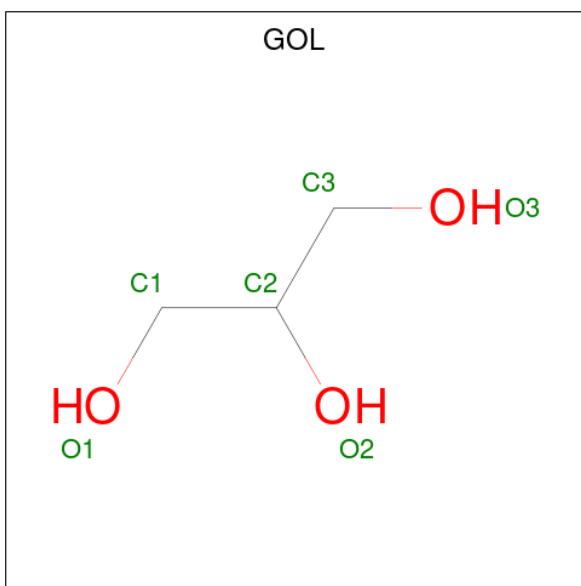
- Molecule 4 is 2-(N-MORPHOLINO)-ETHANESULFONIC ACID (CCD ID: MES) (formula:

C₆H₁₃NO₄S).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	A	1	Total	C	N	O	S	0	0
			12	6	1	4	1		
4	B	1	Total	C	N	O	S	0	0
			12	6	1	4	1		
4	B	1	Total	C	N	O	S	0	0
			12	6	1	4	1		

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



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

- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	196	Total	O	0	0
			196	196		
6	B	153	Total	O	0	0
			153	153		

- Molecule 1: Lin0857 protein



4 Data and refinement statistics

Property	Value	Source
Space group	P 31 2 1	Depositor
Cell constants a, b, c, α , β , γ	145.54Å 145.54Å 105.74Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	43.44 – 2.10 43.44 – 2.10	Depositor EDS
% Data completeness (in resolution range)	99.3 (43.44-2.10) 99.3 (43.44-2.10)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.99 (at 2.10Å)	Xtriage
Refinement program	REFMAC 5.8.0107	Depositor
R, R_{free}	0.179 , 0.215 0.192 , 0.222	Depositor DCC
R_{free} test set	3783 reflections (5.01%)	wwPDB-VP
Wilson B-factor (Å ²)	33.2	Xtriage
Anisotropy	0.028	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 36.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.52$, $\langle L^2 \rangle = 0.36$	Xtriage
Estimated twinning fraction	0.014 for -h,-k,l	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	6149	wwPDB-VP
Average B, all atoms (Å ²)	39.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.57% 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: MAN, MES, SO4, 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	1.34	12/2883 (0.4%)	1.13	3/3910 (0.1%)
1	B	1.37	10/2892 (0.3%)	1.10	2/3921 (0.1%)
All	All	1.35	22/5775 (0.4%)	1.11	5/7831 (0.1%)

The worst 5 of 22 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	276	GLY	N-CA	8.65	1.52	1.44
1	B	138	GLY	C-O	7.52	1.29	1.23
1	B	265	HIS	CA-C	-6.69	1.44	1.52
1	B	11	LEU	C-O	-6.22	1.16	1.24
1	B	4	TYR	CA-C	6.01	1.60	1.53

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	138	GLY	N-CA-C	6.68	118.93	111.85
1	A	3	ILE	N-CA-C	6.26	116.40	106.88
1	A	138	GLY	N-CA-C	5.72	117.91	111.85
1	A	52	VAL	CB-CA-C	-5.65	103.91	110.91
1	B	352	VAL	N-CA-C	5.35	115.56	110.42

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	2815	0	2730	12	0
1	B	2824	0	2743	13	0
2	C	34	0	30	4	0
2	D	34	0	28	4	0
3	A	25	0	0	0	0
3	B	20	0	0	0	0
4	A	12	0	13	2	0
4	B	24	0	26	0	0
5	B	12	0	16	0	0
6	A	196	0	0	0	0
6	B	153	0	0	5	0
All	All	6149	0	5586	26	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 2.

The worst 5 of 26 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:89:ARG:HH12	2:C:1:MAN:H61	1.51	0.73
2:D:1:MAN:O2	2:D:3:BMA:C1	2.40	0.70
1:B:89:ARG:HH22	2:D:1:MAN:H61	1.56	0.69
1:B:3:ILE:HD12	1:B:341:MET:HE2	1.85	0.58
1:A:20:ILE:HD12	4:A:409:MES:H62	1.86	0.56

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	352/363 (97%)	338 (96%)	14 (4%)	0	100	100
1	B	353/363 (97%)	341 (97%)	12 (3%)	0	100	100
All	All	705/726 (97%)	679 (96%)	26 (4%)	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	302/311 (97%)	300 (99%)	2 (1%)	76	83
1	B	303/311 (97%)	300 (99%)	3 (1%)	68	76
All	All	605/622 (97%)	600 (99%)	5 (1%)	73	81

All (5) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	71	LEU
1	A	106	LEU
1	B	15	LEU
1	B	18	LYS
1	B	106	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	146	GLN
1	A	198	ASN
1	A	271	ASN
1	B	154	ASN
1	B	179	GLN

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	MAN	C	1	2	12,12,12	1.41	3 (25%)	17,17,17	3.37	9 (52%)
2	BMA	C	2	2	11,11,12	1.82	1 (9%)	15,15,17	1.78	3 (20%)
2	BMA	C	3	2	11,11,12	0.81	0	15,15,17	1.81	4 (26%)
2	MAN	D	1	2	12,12,12	1.78	5 (41%)	17,17,17	4.04	10 (58%)
2	BMA	D	2	2	11,11,12	1.06	1 (9%)	15,15,17	1.19	1 (6%)
2	BMA	D	3	2	11,11,12	4.76	4 (36%)	15,15,17	4.91	10 (66%)

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	MAN	C	1	2	-	2/2/22/22	0/1/1/1
2	BMA	C	2	2	-	0/2/19/22	0/1/1/1
2	BMA	C	3	2	-	0/2/19/22	1/1/1/1
2	MAN	D	1	2	-	2/2/22/22	0/1/1/1
2	BMA	D	2	2	-	0/2/19/22	0/1/1/1
2	BMA	D	3	2	-	0/2/19/22	1/1/1/1

The worst 5 of 14 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	3	BMA	C1-C2	-10.98	1.25	1.52
2	D	3	BMA	O2-C2	8.16	1.60	1.43
2	D	3	BMA	C2-C3	7.33	1.63	1.52
2	C	2	BMA	O2-C2	-5.43	1.32	1.43
2	D	1	MAN	C1-C2	2.61	1.58	1.52

The worst 5 of 37 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	3	BMA	C1-O5-C5	11.10	127.06	112.19
2	D	3	BMA	C1-C2-C3	8.91	122.61	109.64
2	D	1	MAN	O4-C4-C3	8.49	130.40	110.38
2	D	1	MAN	O2-C2-C1	8.01	127.73	109.25
2	D	3	BMA	O2-C2-C1	7.52	126.45	109.22

There are no chirality outliers.

All (4) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	D	1	MAN	C4-C5-C6-O6
2	C	1	MAN	C4-C5-C6-O6
2	D	1	MAN	O5-C5-C6-O6
2	C	1	MAN	O5-C5-C6-O6

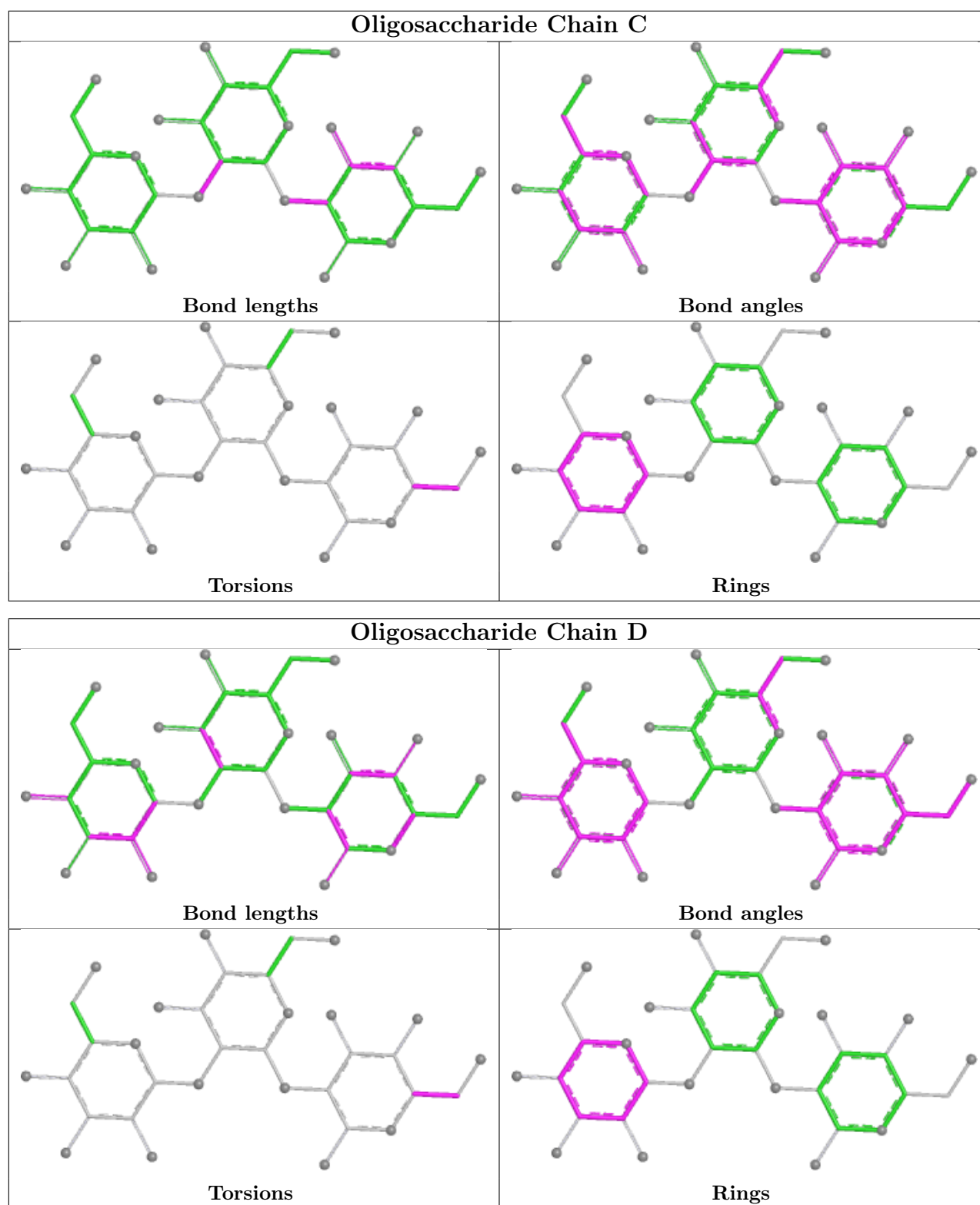
All (2) ring outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	D	3	BMA	C1-C2-C3-C4-C5-O5
2	C	3	BMA	C1-C2-C3-C4-C5-O5

3 monomers are involved in 8 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	D	3	BMA	1	0
2	D	1	MAN	4	0
2	C	1	MAN	4	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](#)

14 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
4	MES	A	409	-	12,12,12	1.94	4 (33%)	15,16,16	3.68	8 (53%)
3	SO4	A	408	-	4,4,4	0.66	0	6,6,6	0.59	0
3	SO4	B	402	-	4,4,4	0.32	0	6,6,6	0.56	0
3	SO4	B	403	-	4,4,4	0.66	0	6,6,6	0.40	0
3	SO4	A	404	-	4,4,4	0.36	0	6,6,6	0.44	0
3	SO4	A	407	-	4,4,4	0.78	0	6,6,6	0.45	0
4	MES	B	407	-	12,12,12	2.11	3 (25%)	15,16,16	2.35	5 (33%)
3	SO4	B	404	-	4,4,4	1.14	0	6,6,6	0.39	0
5	GOL	B	406	-	5,5,5	0.50	0	5,5,5	0.81	0
3	SO4	B	401	-	4,4,4	0.56	0	6,6,6	0.95	0
4	MES	B	408	-	12,12,12	1.79	3 (25%)	15,16,16	3.81	5 (33%)
3	SO4	A	405	-	4,4,4	0.69	0	6,6,6	0.54	0
5	GOL	B	405	-	5,5,5	0.65	0	5,5,5	0.72	0
3	SO4	A	406	-	4,4,4	0.32	0	6,6,6	0.42	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	MES	A	409	-	-	5/6/14/14	0/1/1/1
4	MES	B	407	-	-	2/6/14/14	0/1/1/1
5	GOL	B	406	-	-	3/4/4/4	-
4	MES	B	408	-	-	0/6/14/14	0/1/1/1
5	GOL	B	405	-	-	0/4/4/4	-

The worst 5 of 10 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	B	407	MES	C8-S	-5.04	1.70	1.77
4	A	409	MES	O2S-S	3.49	1.55	1.45
4	B	407	MES	O2S-S	3.40	1.54	1.45
4	B	408	MES	O1S-S	3.39	1.54	1.45

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	B	408	MES	C8-S	-3.28	1.73	1.77

The worst 5 of 18 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	408	MES	O2S-S-C8	-9.59	92.23	106.73
4	B	408	MES	O3S-S-C8	9.49	124.57	106.00
4	A	409	MES	O1S-S-C8	8.36	119.36	106.73
4	A	409	MES	O3S-S-O2S	6.17	126.82	111.40
4	B	407	MES	O1S-S-C8	5.53	115.09	106.73

There are no chirality outliers.

5 of 10 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	409	MES	C8-C7-N4-C5
4	A	409	MES	C7-C8-S-O2S
5	B	406	GOL	O1-C1-C2-O2
4	A	409	MES	C7-C8-S-O3S
5	B	406	GOL	O1-C1-C2-C3

There are no ring outliers.

1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	409	MES	2	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 [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	354/363 (97%)	-0.01	8 (2%) 61 64	26, 36, 52, 72	1 (0%)
1	B	355/363 (97%)	0.15	7 (1%) 65 67	27, 39, 58, 84	1 (0%)
All	All	709/726 (97%)	0.07	15 (2%) 63 66	26, 37, 56, 84	2 (0%)

The worst 5 of 15 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	36	GLU	4.2
1	B	257	GLU	3.3
1	A	23	GLY	3.2
1	A	36	GLU	2.9
1	B	355	LYS	2.9

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.

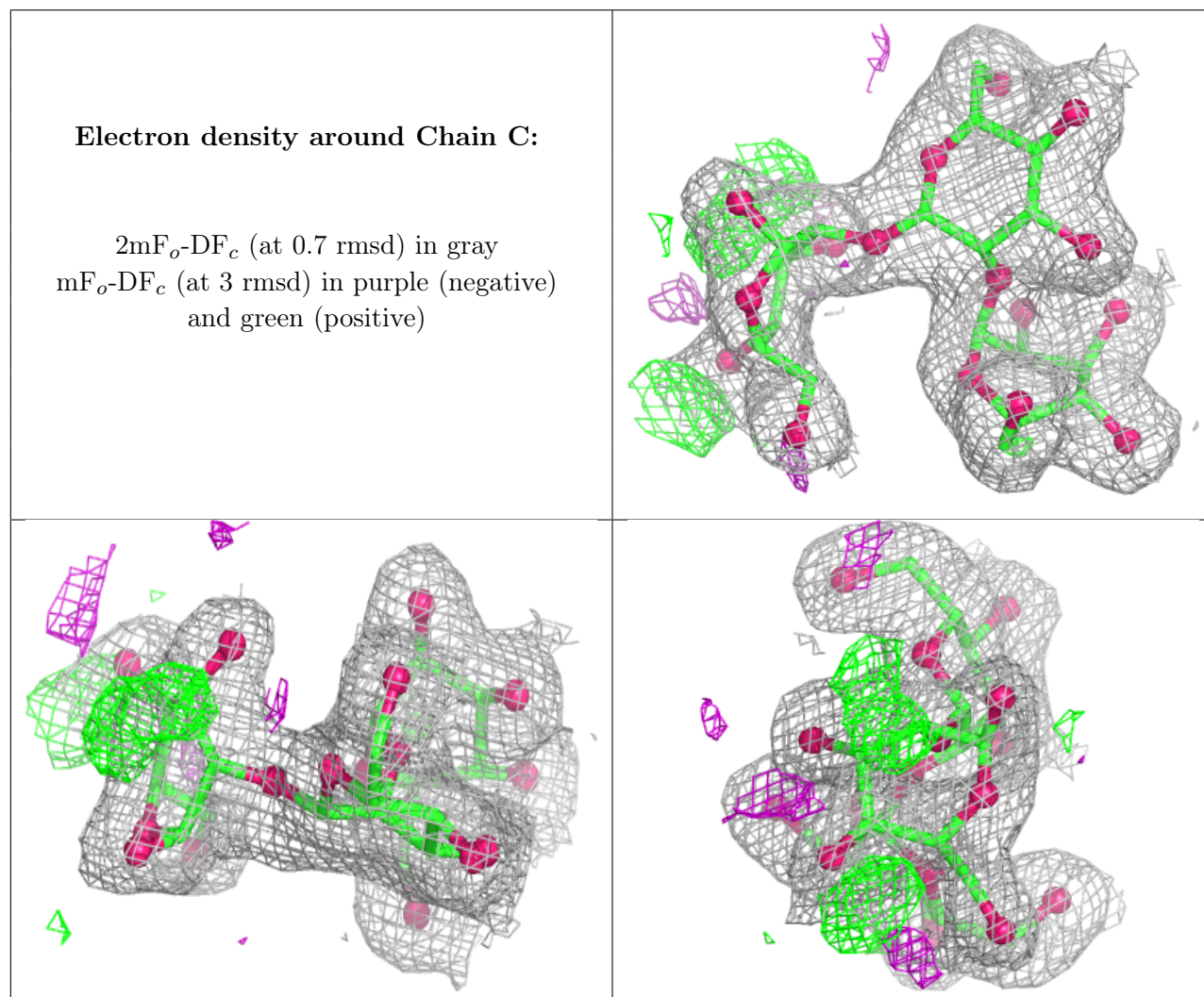
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	MAN	D	1	12/12	0.78	0.21	34,50,65,66	0
2	MAN	C	1	12/12	0.88	0.15	33,49,61,63	0
2	BMA	D	2	11/12	0.95	0.07	27,29,31,32	0
2	BMA	D	3	11/12	0.95	0.08	25,28,30,32	0
2	BMA	C	3	11/12	0.98	0.06	24,26,29,29	0

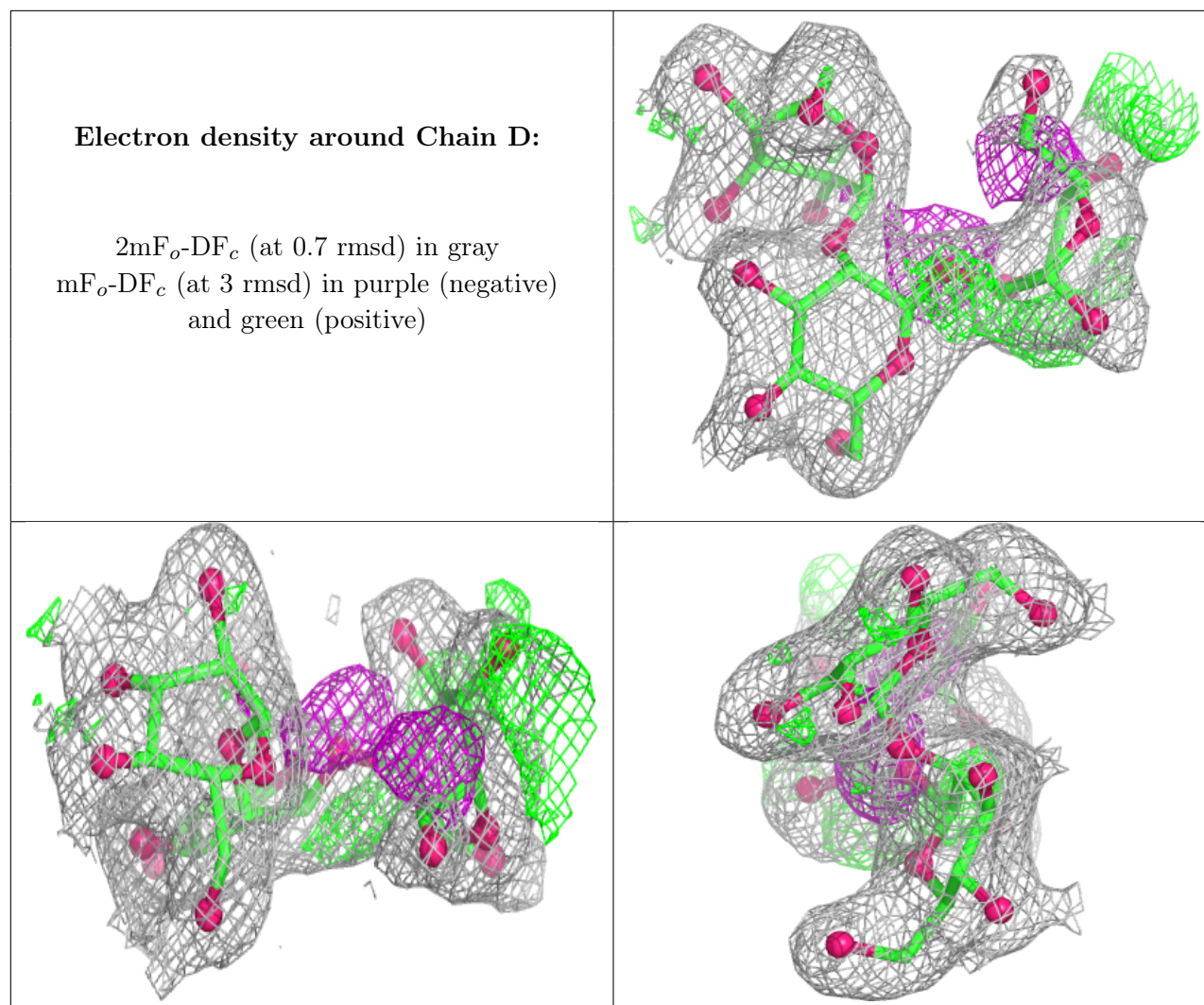
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	BMA	C	2	11/12	0.98	0.05	26,27,29,30	0

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 ⓘ

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	SO4	B	404	5/5	0.72	0.14	68,74,80,86	0
3	SO4	B	403	5/5	0.73	0.12	76,78,84,91	0
4	MES	A	409	12/12	0.73	0.22	55,58,62,63	0
3	SO4	A	406	5/5	0.78	0.16	92,95,101,108	0
3	SO4	A	408	5/5	0.81	0.12	61,63,80,83	0
5	GOL	B	406	6/6	0.88	0.17	56,64,67,68	0
4	MES	B	408	12/12	0.90	0.15	49,51,58,62	0
3	SO4	A	405	5/5	0.91	0.09	54,55,71,76	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
5	GOL	B	405	6/6	0.93	0.11	39,53,54,65	0
4	MES	B	407	12/12	0.93	0.13	46,48,52,53	0
3	SO4	B	401	5/5	0.94	0.08	59,59,64,64	0
3	SO4	A	407	5/5	0.95	0.17	54,54,57,66	0
3	SO4	B	402	5/5	0.99	0.04	32,33,34,34	0
3	SO4	A	404	5/5	1.00	0.03	28,28,29,30	0

6.5 Other polymers [i](#)

There are no such residues in this entry.