



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 10, 2026 – 05:10 PM UTC

PDB ID : 5AZA / pdb_00005aza
Title : Crystal structure of MBP-sAglB fusion protein with a 20-residue spacer in the connector helix
Authors : Matsuoka, R.; Kohda, D.
Deposited on : 2015-09-27
Resolution : 2.08 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

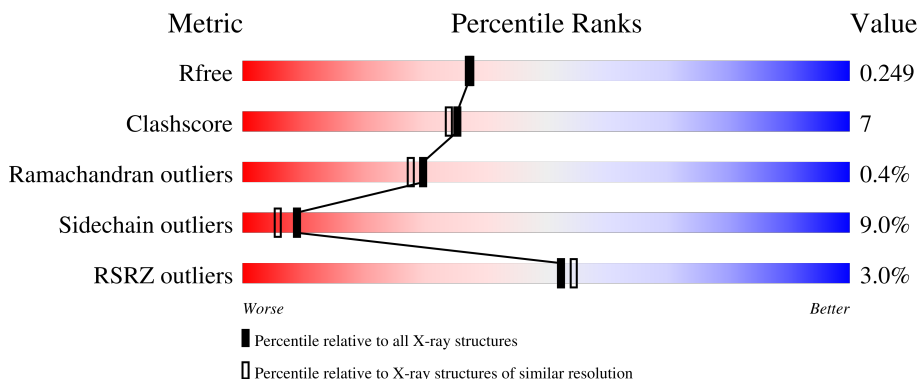
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

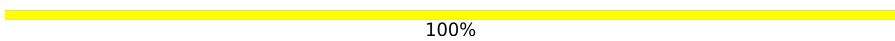
The reported resolution of this entry is 2.08 Å.

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	8172 (2.10-2.06)
Clashscore	190562	8714 (2.10-2.06)
Ramachandran outliers	187476	8641 (2.10-2.06)
Sidechain outliers	187428	8642 (2.10-2.06)
RSRZ outliers	180081	8177 (2.10-2.06)

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	872	
2	B	2	

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 6733 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 Maltose-binding periplasmic protein,Oligosaccharyl transferase stt3 subunit related protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	836	Total	C	N	O	S	0	4	0
			6583	4239	1076	1256	12			

There are 28 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	0	MET	-	initiating methionine	UNP P0AEX9
A	312	VAL	ALA	engineered mutation	UNP P0AEX9
A	369	ALA	-	linker	UNP P0AEX9
A	370	ALA	-	linker	UNP P0AEX9
A	371	ALA	-	linker	UNP P0AEX9
A	372	LYS	-	linker	UNP P0AEX9
A	373	ALA	-	linker	UNP P0AEX9
A	374	GLU	-	linker	UNP P0AEX9
A	375	ALA	-	linker	UNP P0AEX9
A	376	ALA	-	linker	UNP P0AEX9
A	377	ALA	-	linker	UNP P0AEX9
A	378	LYS	-	linker	UNP P0AEX9
A	379	ALA	-	linker	UNP P0AEX9
A	380	GLU	-	linker	UNP P0AEX9
A	381	ALA	-	linker	UNP P0AEX9
A	382	ALA	-	linker	UNP P0AEX9
A	383	ALA	-	linker	UNP P0AEX9
A	384	LYS	-	linker	UNP P0AEX9
A	385	ALA	-	linker	UNP P0AEX9
A	386	ALA	-	linker	UNP P0AEX9
A	387	ALA	-	linker	UNP P0AEX9
A	388	ALA	-	linker	UNP P0AEX9
A	866	HIS	-	expression tag	UNP I6V0B8
A	867	HIS	-	expression tag	UNP I6V0B8
A	868	HIS	-	expression tag	UNP I6V0B8
A	869	HIS	-	expression tag	UNP I6V0B8

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Chain	Residue	Modelled	Actual	Comment	Reference
A	870	HIS	-	expression tag	UNP I6V0B8
A	871	HIS	-	expression tag	UNP I6V0B8

- Molecule 2 is an oligosaccharide called alpha-D-glucopyranose-(1-4)-alpha-D-glucopyranose.



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace
2	B	2	Total	C	O	0	0	0
			23	12	11			

- Molecule 3 is CALCIUM ION (CCD ID: CA) (formula: Ca).

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

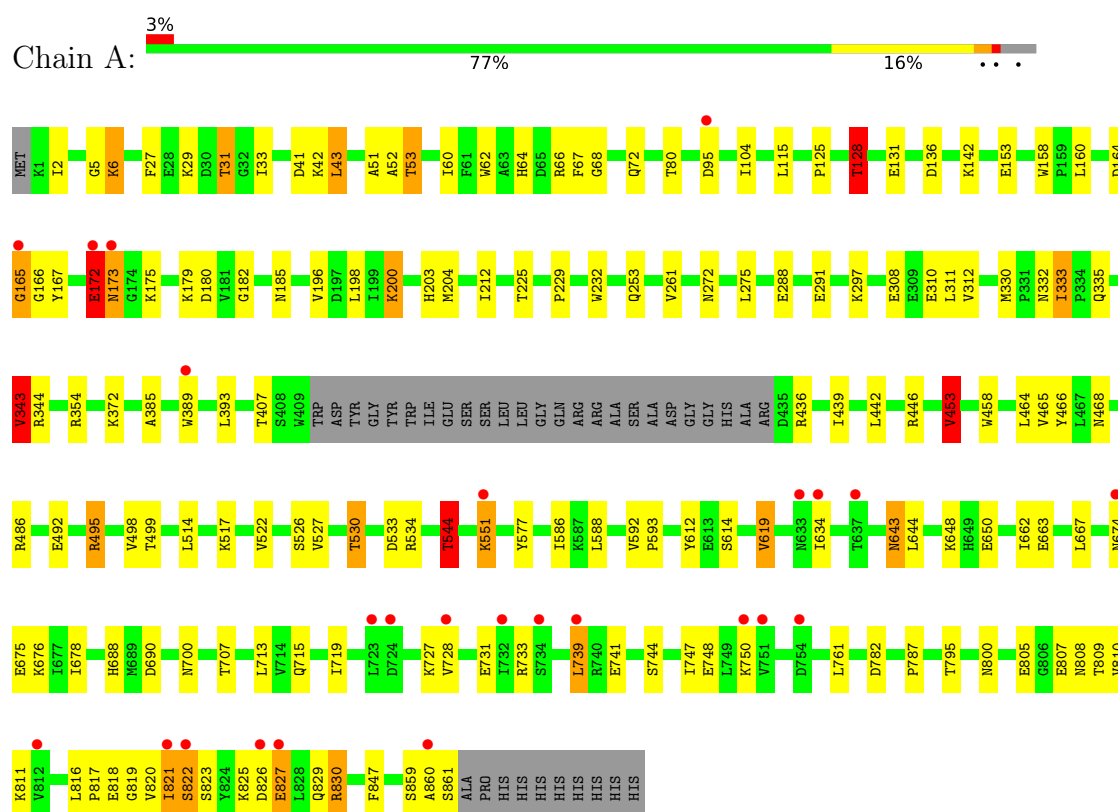
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	126	Total	O	0	0
			126	126		

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: Maltose-binding periplasmic protein,Oligosaccharyl transferase stt3 subunit related protein



- Molecule 2: alpha-D-glucopyranose-(1-4)-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	66.47Å 100.47Å 140.92Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	50.00 – 2.08 50.00 – 2.08	Depositor EDS
% Data completeness (in resolution range)	99.7 (50.00-2.08) 99.8 (50.00-2.08)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.33 (at 2.08Å)	Xtriage
Refinement program	REFMAC 5.8.0073	Depositor
R, R_{free}	0.196 , 0.248 0.204 , 0.249	Depositor DCC
R_{free} test set	2914 reflections (5.05%)	wwPDB-VP
Wilson B-factor (Å ²)	34.5	Xtriage
Anisotropy	0.026	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 29.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	6733	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 4.33% 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: GLC, CA

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	1.13	3/6733 (0.0%)	1.11	14/9140 (0.2%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	2

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	172	GLU	CA-C	-6.30	1.47	1.52
1	A	787	PRO	C-O	-5.64	1.17	1.23
1	A	225	THR	CA-C	5.02	1.59	1.52

All (14) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	53	THR	N-CA-C	-12.11	90.90	109.79
1	A	821	ILE	N-CA-C	-8.13	101.18	110.05
1	A	128	THR	N-CA-CB	-6.48	99.55	110.83
1	A	53	THR	CA-C-N	-6.39	108.89	121.41
1	A	53	THR	C-N-CA	-6.39	108.89	121.41
1	A	544	THR	CB-CA-C	5.76	117.52	108.88
1	A	527	VAL	CB-CA-C	-5.46	102.36	110.33
1	A	53	THR	O-C-N	-5.41	114.66	121.74
1	A	453	VAL	N-CA-CB	-5.29	100.16	111.37
1	A	343	VAL	CB-CA-C	5.26	120.73	112.16

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	80	THR	CB-CA-C	-5.18	104.26	111.41
1	A	125	PRO	CA-C-N	5.09	125.00	119.76
1	A	125	PRO	C-N-CA	5.09	125.00	119.76
1	A	830	ARG	N-CA-C	-5.05	105.46	110.97

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	166	GLY	Peptide
1	A	172	GLU	Peptide

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	6583	0	6540	98	0
2	B	23	0	21	0	0
3	A	1	0	0	0	0
4	A	126	0	0	13	0
All	All	6733	0	6561	98	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 7.

All (98) 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:95:ASP:HB3	4:A:1115:HOH:O	1.26	1.26
1:A:522:VAL:HB	4:A:1091:HOH:O	1.52	1.08
1:A:167:TYR:O	1:A:179:LYS:HE3	1.73	0.89
1:A:530:THR:HG22	4:A:1046:HOH:O	1.74	0.88
1:A:72:GLN:HG3	4:A:1005:HOH:O	1.80	0.82
1:A:203:HIS:HE1	4:A:1118:HOH:O	1.64	0.79
1:A:64:HIS:HD2	1:A:261:VAL:H	1.31	0.75

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:688:HIS:HE1	1:A:690:ASP:OD1	1.72	0.71
1:A:522:VAL:HG12	1:A:522:VAL:O	1.91	0.70
1:A:128:THR:HB	1:A:131:GLU:OE1	1.92	0.68
1:A:663:GLU:OE2	1:A:688:HIS:HD2	1.76	0.67
1:A:64:HIS:CD2	1:A:261:VAL:H	2.12	0.67
1:A:643:ASN:HD22	1:A:643:ASN:C	2.03	0.67
1:A:827:GLU:HA	1:A:830:ARG:HG3	1.79	0.65
1:A:819:GLY:O	1:A:823:SER:CB	2.45	0.64
1:A:486:ARG:NH1	1:A:782:ASP:OD1	2.27	0.64
1:A:27:PHE:O	1:A:31:THR:HB	1.98	0.63
1:A:196:VAL:CG1	1:A:200:LYS:HD2	2.29	0.63
1:A:95:ASP:CB	4:A:1115:HOH:O	2.07	0.62
1:A:308:GLU:O	1:A:312:VAL:HG23	2.00	0.62
1:A:182:GLY:O	1:A:185:ASN:ND2	2.31	0.62
1:A:544:THR:CG2	4:A:1077:HOH:O	2.45	0.62
1:A:196:VAL:HG12	1:A:200:LYS:HD2	1.80	0.61
1:A:644:LEU:O	1:A:860:ALA:HA	2.01	0.60
1:A:31:THR:HG22	1:A:33:ILE:H	1.67	0.59
1:A:747:ILE:O	1:A:811:LYS:HA	2.01	0.59
1:A:819:GLY:O	1:A:823:SER:HB3	2.02	0.59
1:A:164:ASP:O	1:A:165:GLY:O	2.21	0.58
1:A:820:VAL:HG23	1:A:821:ILE:H	1.70	0.57
1:A:198:LEU:HD13	1:A:204:MET:HE3	1.87	0.56
1:A:544:THR:HG23	4:A:1077:HOH:O	2.06	0.56
1:A:196:VAL:HG12	1:A:200:LYS:CD	2.35	0.56
1:A:739:LEU:HD11	1:A:747:ILE:HD11	1.86	0.56
1:A:68:GLY:HA3	1:A:332:ASN:O	2.06	0.56
1:A:297[A]:LYS:HD3	4:A:1052:HOH:O	2.05	0.56
1:A:128:THR:HG22	1:A:131:GLU:H	1.71	0.55
1:A:663:GLU:CD	1:A:688:HIS:HD2	2.14	0.55
1:A:5:GLY:H	1:A:272:ASN:HD21	1.54	0.55
1:A:741:GLU:HA	1:A:816:LEU:HD23	1.88	0.54
1:A:650:GLU:HA	1:A:700:ASN:HD22	1.71	0.54
1:A:136:ASP:OD2	1:A:203:HIS:HD2	1.90	0.54
1:A:51:ALA:O	1:A:53:THR:O	2.26	0.54
1:A:663:GLU:H	1:A:715:GLN:HE22	1.54	0.53
1:A:312:VAL:HG12	1:A:312:VAL:O	2.08	0.53
1:A:164:ASP:C	1:A:165:GLY:O	2.51	0.53
1:A:663:GLU:H	1:A:715:GLN:NE2	2.07	0.53
1:A:727:LYS:HA	1:A:731:GLU:O	2.08	0.53
1:A:822:SER:HA	1:A:825:LYS:HB3	1.91	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:52:ALA:C	1:A:53:THR:O	2.49	0.51
1:A:498:VAL:HG11	1:A:577:TYR:CD1	2.45	0.51
1:A:662:ILE:HA	1:A:715:GLN:HE22	1.76	0.51
1:A:464:LEU:HD12	1:A:464:LEU:N	2.25	0.51
1:A:229:PRO:HA	1:A:232:TRP:CE2	2.46	0.51
1:A:498:VAL:HG11	1:A:577:TYR:CG	2.47	0.50
1:A:64:HIS:HE1	1:A:330:MET:O	1.94	0.49
1:A:643:ASN:OD1	1:A:861:SER:C	2.56	0.49
1:A:33:ILE:HD13	1:A:275:LEU:HD22	1.95	0.48
1:A:31:THR:CG2	1:A:33:ILE:H	2.27	0.48
1:A:643:ASN:HB2	1:A:859:SER:O	2.13	0.48
1:A:67:PHE:HB3	1:A:104:ILE:HD12	1.96	0.47
1:A:544:THR:HG21	1:A:593:PRO:HD2	1.97	0.47
1:A:741:GLU:OE1	1:A:821:ILE:O	2.33	0.47
1:A:690:ASP:HB3	4:A:1088:HOH:O	2.14	0.47
1:A:819:GLY:O	1:A:823:SER:HB2	2.14	0.47
1:A:62:TRP:CD1	1:A:66:ARG:HG3	2.50	0.47
1:A:172:GLU:O	1:A:173:ASN:HB2	2.14	0.47
1:A:612:TYR:HB3	1:A:619:VAL:HG13	1.96	0.47
1:A:62:TRP:HB3	1:A:67:PHE:HE1	1.80	0.46
1:A:820:VAL:HG23	1:A:821:ILE:N	2.29	0.46
1:A:728:VAL:HG23	1:A:733:ARG:HG3	1.98	0.46
1:A:495:ARG:HH11	1:A:495:ARG:HG3	1.81	0.45
1:A:551:LYS:CE	4:A:1016:HOH:O	2.64	0.45
1:A:153:GLU:OE1	1:A:344:ARG:NH1	2.50	0.45
1:A:522:VAL:O	1:A:522:VAL:CG1	2.63	0.45
1:A:31:THR:HG23	1:A:33:ILE:HD12	1.98	0.44
1:A:522:VAL:HA	1:A:526:SER:O	2.17	0.44
1:A:662:ILE:C	1:A:663:GLU:HG2	2.42	0.44
1:A:544:THR:HG22	4:A:1077:HOH:O	2.14	0.44
1:A:442:LEU:HD12	1:A:446:ARG:HG3	2.00	0.44
1:A:466:TYR:CE2	1:A:468:ASN:HB2	2.53	0.43
1:A:385:ALA:O	1:A:389:TRP:HD1	2.01	0.43
1:A:6:LYS:HA	1:A:33:ILE:HG23	2.00	0.43
1:A:43:LEU:C	1:A:43:LEU:HD12	2.44	0.43
1:A:498:VAL:CG1	1:A:499:THR:N	2.82	0.42
1:A:436:ARG:HG2	1:A:458:TRP:CZ2	2.55	0.42
1:A:439:ILE:HG23	1:A:453:VAL:HG22	2.01	0.42
1:A:748:GLU:HA	1:A:810:VAL:O	2.20	0.42
1:A:158:TRP:CH2	1:A:343:VAL:HG22	2.54	0.42
1:A:817:PRO:O	1:A:820:VAL:HG22	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:586:ILE:HG22	1:A:592:VAL:HB	2.02	0.41
1:A:333:ILE:HD12	1:A:335:GLN:HB2	2.03	0.41
1:A:807:GLU:C	1:A:808:ASN:ND2	2.79	0.41
1:A:495:ARG:CZ	1:A:495:ARG:HB3	2.49	0.41
1:A:172:GLU:OE2	1:A:372:LYS:NZ	2.51	0.41
1:A:498:VAL:HG12	1:A:499:THR:N	2.35	0.41
1:A:203:HIS:CE1	4:A:1118:HOH:O	2.52	0.40
1:A:33:ILE:CD1	1:A:275:LEU:HD22	2.51	0.40
1:A:761:LEU:HD13	1:A:847:PHE:CZ	2.57	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	836/872 (96%)	809 (97%)	24 (3%)	3 (0%)	30 28

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	165	GLY
1	A	800	ASN
1	A	674	ASN

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	690/716 (96%)	628 (91%)	62 (9%)	9 6

All (62) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	2	ILE
1	A	6	LYS
1	A	29	LYS
1	A	31	THR
1	A	41	ASP
1	A	42	LYS
1	A	43	LEU
1	A	60	ILE
1	A	115	LEU
1	A	128	THR
1	A	142	LYS
1	A	160	LEU
1	A	173	ASN
1	A	175	LYS
1	A	180	ASP
1	A	200	LYS
1	A	212	ILE
1	A	253	GLN
1	A	288	GLU
1	A	291	GLU
1	A	310	GLU
1	A	311	LEU
1	A	333	ILE
1	A	343	VAL
1	A	354	ARG
1	A	393	LEU
1	A	407	THR
1	A	453	VAL
1	A	465	VAL
1	A	492	GLU
1	A	495	ARG
1	A	514	LEU
1	A	517	LYS
1	A	530	THR
1	A	533	ASP
1	A	534	ARG
1	A	544	THR
1	A	551	LYS

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Mol	Chain	Res	Type
1	A	588	LEU
1	A	614	SER
1	A	619	VAL
1	A	634	ILE
1	A	643	ASN
1	A	648	LYS
1	A	667	LEU
1	A	675	GLU
1	A	676	LYS
1	A	678	ILE
1	A	707	THR
1	A	713	LEU
1	A	719	ILE
1	A	739	LEU
1	A	744	SER
1	A	750	LYS
1	A	795	THR
1	A	805	GLU
1	A	809	THR
1	A	818	GLU
1	A	822	SER
1	A	826	ASP
1	A	827	GLU
1	A	829	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (21) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	49	GLN
1	A	64	HIS
1	A	100	ASN
1	A	173	ASN
1	A	201	ASN
1	A	203	HIS
1	A	218	ASN
1	A	272	ASN
1	A	325	GLN
1	A	399	ASN
1	A	474	ASN
1	A	524	ASN
1	A	607	ASN
1	A	616	ASN

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Mol	Chain	Res	Type
1	A	643	ASN
1	A	673	ASN
1	A	688	HIS
1	A	700	ASN
1	A	715	GLN
1	A	737	ASN
1	A	829	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 ⓘ

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
2	GLC	B	1	2	12,12,12	0.74	0	17,17,17	1.55	3 (17%)
2	GLC	B	2	2	11,11,12	0.51	0	15,15,17	1.35	2 (13%)

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	GLC	B	1	2	-	0/2/22/22	0/1/1/1
2	GLC	B	2	2	-	0/2/19/22	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(°)
2	B	1	GLC	O5-C1-C2	3.82	117.02	110.30
2	B	2	GLC	C1-O5-C5	2.82	115.96	112.19
2	B	2	GLC	C2-C3-C4	-2.63	106.24	110.86
2	B	1	GLC	O2-C2-C3	-2.13	105.35	110.38
2	B	1	GLC	O4-C4-C3	-2.04	105.56	110.38

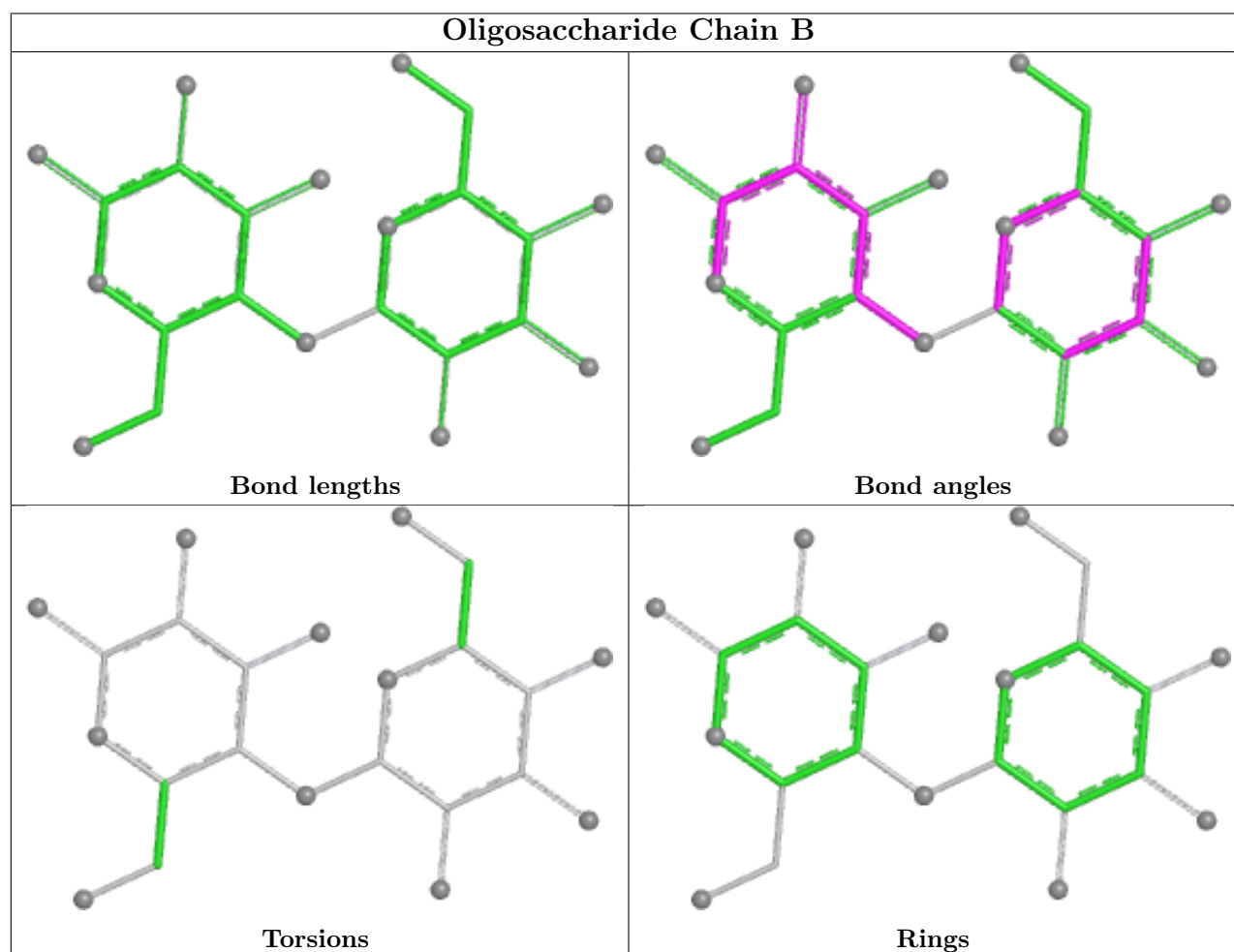
There are no chirality outliers.

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

Of 1 ligands modelled in this entry, 1 is monoatomic - leaving 0 for Mogul analysis.

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.

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, 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	836/872 (95%)	0.17	25 (2%)	52 54	13, 35, 67, 91	4 (0%)

All (25) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	165	GLY	4.4
1	A	95	ASP	3.8
1	A	173	ASN	3.3
1	A	634	ILE	3.2
1	A	812	VAL	3.0
1	A	822	SER	2.9
1	A	739	LEU	2.8
1	A	637	THR	2.6
1	A	389	TRP	2.6
1	A	723	LEU	2.6
1	A	633	ASN	2.3
1	A	732	ILE	2.3
1	A	860	ALA	2.3
1	A	827	GLU	2.2
1	A	724	ASP	2.2
1	A	734	SER	2.2
1	A	728	VAL	2.2
1	A	172	GLU	2.1
1	A	821	ILE	2.1
1	A	674	ASN	2.1
1	A	750	LYS	2.1
1	A	751	VAL	2.1
1	A	826	ASP	2.0
1	A	551	LYS	2.0
1	A	754	ASP	2.0

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

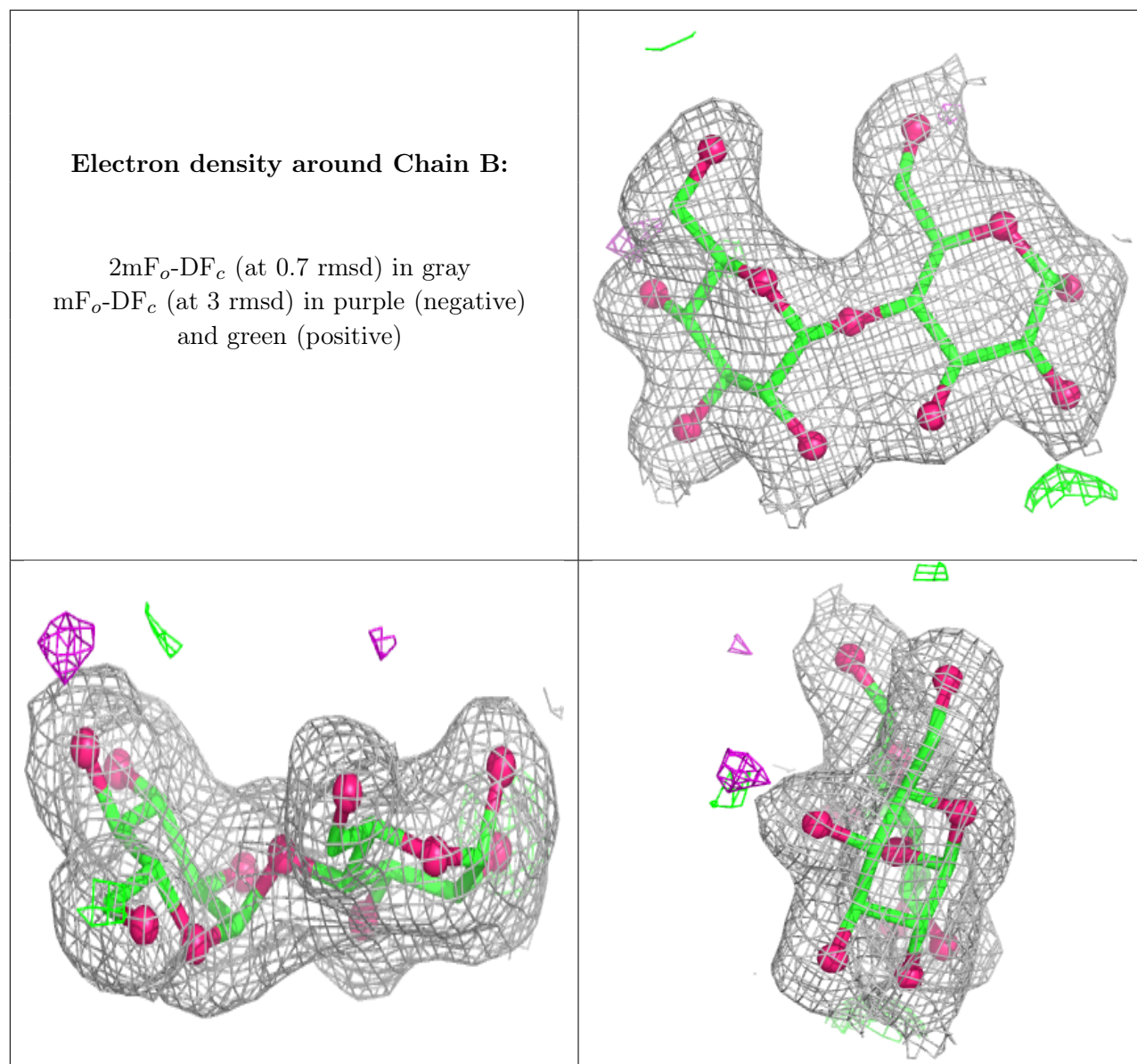
There are no non-standard protein/DNA/RNA residues in this entry.

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($\text{\AA}^2$)	Q<0.9
2	GLC	B	2	11/12	0.95	0.06	22,24,26,27	0
2	GLC	B	1	12/12	0.97	0.04	23,27,31,31	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(Å ²)	Q<0.9
3	CA	A	902	1/1	0.98	0.03	25,25,25,25	0

6.5 Other polymers [i](#)

There are no such residues in this entry.