



# Full wwPDB X-ray Structure Validation Report ⓘ

Sep 3, 2025 – 06:04 PM EDT

PDB ID : 9PX5 / pdb\_00009px5  
Title : Crystal structure of Fab 7268 in complex with MBP-TREM2 Ig domain fusion  
Authors : Arndt, J.W.; Chao, Q.; Cooke, H.A.; Almeida, A.D.S.  
Deposited on : 2025-08-05  
Resolution : 3.70 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 : 4-5-2 with Phenix2.0rc1  
Mogul : 2022.3.0, CSD as543be (2022)  
Xtriage (Phenix) : 2.0rc1  
EDS : 3.0  
Percentile statistics : 20231227.v01 (using entries in the PDB archive December 27th 2023)  
CCP4 : 9.0.006 (Gargrove)  
Density-Fitness : 1.0.12  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.45.1

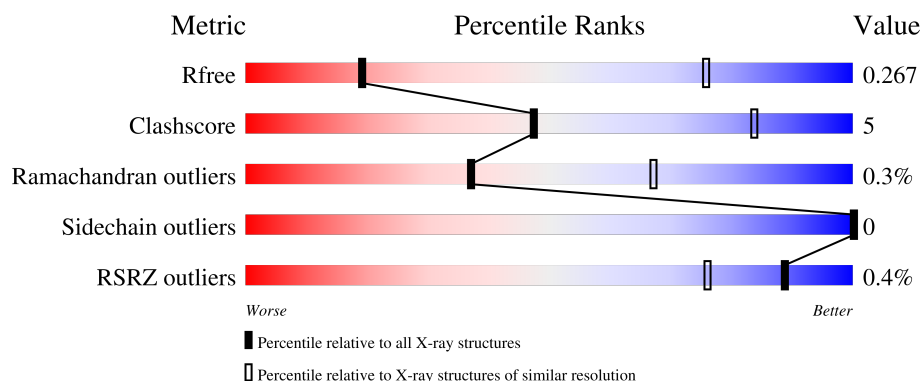
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*





The reported resolution of this entry is 3.70 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	1017 (3.80-3.60)
Clashscore	180529	1074 (3.80-3.60)
Ramachandran outliers	177936	1055 (3.80-3.60)
Sidechain outliers	177891	1052 (3.80-3.60)
RSRZ outliers	164620	1017 (3.80-3.60)

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 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	A	492	
2	H	226	
3	L	214	
4	B	2	

## 2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 6597 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/maltodextrin-binding periplasmic protein, Triggering receptor expressed on myeloid cells 2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	464	Total	C	N	O	S	0	0	0
			3324	2115	549	649	11			

There are 22 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-348	THR	ILE	conflict	UNP P0AEX9
A	-268	ALA	ASP	conflict	UNP P0AEX9
A	-267	ALA	LYS	conflict	UNP P0AEX9
A	-178	ALA	GLU	conflict	UNP P0AEX9
A	-177	ALA	ASN	conflict	UNP P0AEX9
A	-135	HIS	ALA	conflict	UNP P0AEX9
A	-131	HIS	LYS	conflict	UNP P0AEX9
A	-111	ALA	LYS	conflict	UNP P0AEX9
A	-38	VAL	ALA	conflict	UNP P0AEX9
A	-33	VAL	ILE	conflict	UNP P0AEX9
A	9	ALA	GLU	conflict	UNP P0AEX9
A	12	ALA	LYS	conflict	UNP P0AEX9
A	13	ALA	ASP	conflict	UNP P0AEX9
A	17	ASN	-	linker	UNP P0AEX9
A	20	ASP	ASN	conflict	UNP Q9NZC2
A	136	HIS	-	expression tag	UNP Q9NZC2
A	137	HIS	-	expression tag	UNP Q9NZC2
A	138	HIS	-	expression tag	UNP Q9NZC2
A	139	HIS	-	expression tag	UNP Q9NZC2
A	140	HIS	-	expression tag	UNP Q9NZC2
A	141	HIS	-	expression tag	UNP Q9NZC2
A	142	HIS	-	expression tag	UNP Q9NZC2

- Molecule 2 is a protein called 7268 Fab Heavy Chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	H	225	Total	C	N	O	S	0	0	0
			1642	1043	267	327	5			

- Molecule 3 is a protein called 7268 Fab Light Chain.

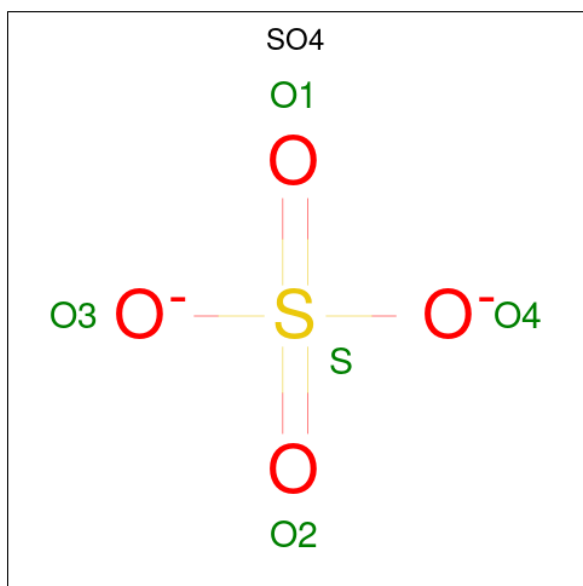
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	L	214	Total	C	N	O	S	0	0	0
			1598	997	263	333	5			

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



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

- Molecule 5 is SULFATE ION (CCD ID: SO4) (formula: O<sub>4</sub>S).

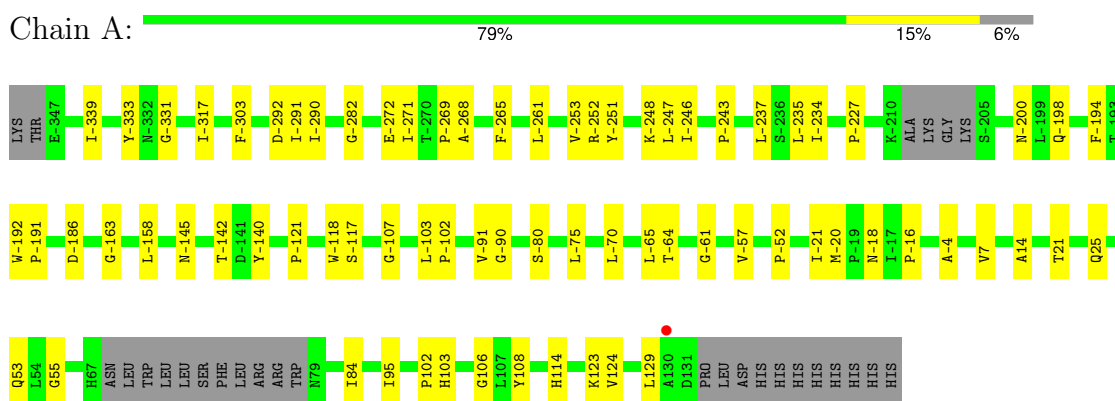


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	O	S	0	0
			5	4	1		
5	A	1	Total	O	S	0	0
			5	4	1		

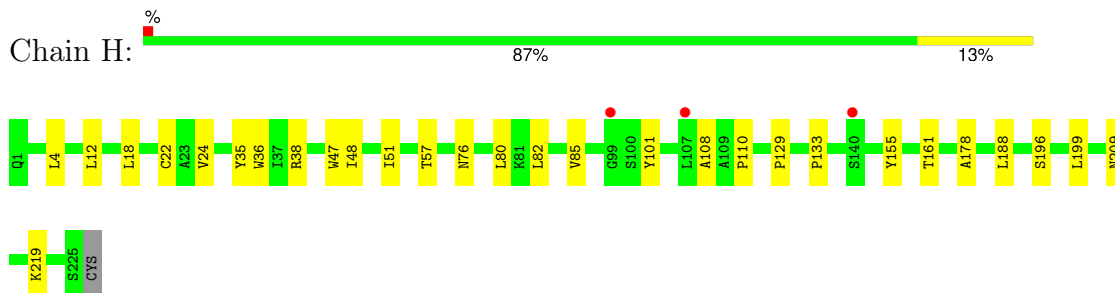
### 3 Residue-property plots

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.

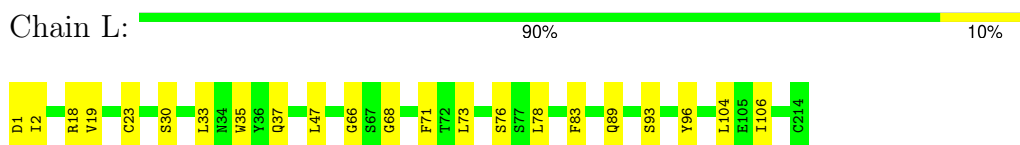
- Molecule 1: Maltose/maltodextrin-binding periplasmic protein, Triggering receptor expressed on myeloid cells 2



- Molecule 2: 7268 Fab Heavy Chain



- Molecule 3: 7268 Fab Light Chain



- Molecule 4: alpha-D-glucopyranose-(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, $\alpha$ , $\beta$ , $\gamma$	119.90Å 64.23Å 160.80Å 90.00° 94.45° 90.00°	Depositor
Resolution (Å)	47.05 – 3.70 47.05 – 3.70	Depositor EDS
% Data completeness (in resolution range)	99.2 (47.05-3.70) 99.1 (47.05-3.70)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.36 (at 3.57Å)	Xtriage
Refinement program	PHENIX 1.21.2-5419_1692	Depositor
R, $R_{free}$	0.248 , 0.268 0.247 , 0.267	Depositor DCC
$R_{free}$ test set	719 reflections (4.98%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	116.4	Xtriage
Anisotropy	0.202	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.30 , 81.5	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.47$ , $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.90	EDS
Total number of atoms	6597	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	123.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.07% 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: GLC, SO4, BGC

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	0.09	0/3406	0.26	0/4687
2	H	0.10	0/1689	0.31	0/2323
3	L	0.08	0/1633	0.27	0/2229
All	All	0.09	0/6728	0.28	0/9239

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	3324	0	2978	42	0
2	H	1642	0	1539	15	0
3	L	1598	0	1493	12	0
4	B	23	0	21	0	0
5	A	10	0	0	0	0
All	All	6597	0	6031	69	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 5.

All (69) 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:102:PRO:O	1:A:103:HIS:ND1	2.23	0.71
1:A:-64:THR:HG23	1:A:-61:GLY:H	1.56	0.71
1:A:-200:ASN:HD22	1:A:-140:TYR:HB2	1.55	0.70
2:H:129:PRO:HB3	2:H:155:TYR:HB3	1.74	0.70
2:H:24:VAL:O	2:H:76:ASN:ND2	2.28	0.67
1:A:-145:ASN:HB2	1:A:-142:THR:HG23	1.85	0.59
2:H:51:ILE:HG22	2:H:57:THR:HG22	1.84	0.59
1:A:-158:LEU:HD23	1:A:7:VAL:HG13	1.85	0.58
2:H:38:ARG:HB3	2:H:48:ILE:HD11	1.86	0.57
2:H:133:PRO:HD3	2:H:219:LYS:HE2	1.86	0.57
1:A:25:GLN:HB2	1:A:129:LEU:HD11	1.87	0.56
1:A:-291:ILE:HD13	1:A:-70:LEU:HD21	1.88	0.55
1:A:-118:TRP:HB2	1:A:-52:PRO:HG2	1.88	0.55
3:L:37:GLN:HB2	3:L:47:LEU:HD11	1.88	0.55
1:A:-282:GLY:HA3	1:A:-18:ASN:O	2.08	0.54
1:A:-91:VAL:HB	1:A:-21:ILE:HA	1.90	0.54
1:A:-235:LEU:HD22	1:A:-102:PRO:HD3	1.92	0.52
1:A:-235:LEU:HB2	1:A:-103:LEU:HD23	1.92	0.51
1:A:-268:ALA:O	1:A:-265:PHE:N	2.43	0.51
1:A:-121:PRO:HA	1:A:-118:TRP:CE2	2.46	0.51
1:A:-117:SER:HB3	1:A:-52:PRO:HD3	1.93	0.51
1:A:-317:ILE:HD13	1:A:-75:LEU:HD13	1.93	0.50
2:H:35:TYR:HD2	2:H:47:TRP:HE1	1.60	0.50
2:H:178:ALA:HA	2:H:188:LEU:HB3	1.93	0.50
1:A:-271:ILE:HG22	1:A:-269:PRO:HD3	1.94	0.50
1:A:-303:PHE:CG	1:A:-290:ILE:HD12	2.46	0.50
3:L:30:SER:O	3:L:68:GLY:N	2.42	0.50
1:A:-251:TYR:HE2	1:A:-16:PRO:HD3	1.77	0.49
1:A:-192:TRP:N	1:A:-191:PRO:HD2	2.28	0.49
2:H:12:LEU:HD21	2:H:18:LEU:HA	1.95	0.49
3:L:1:ASP:OD1	3:L:1:ASP:N	2.44	0.48
3:L:2:ILE:HD12	3:L:93:SER:HB2	1.96	0.48
1:A:-292:ASP:OD2	1:A:-80:SER:OG	2.24	0.48
1:A:-265:PHE:HZ	1:A:-65:LEU:HD12	1.79	0.48
1:A:-261:LEU:HD11	1:A:-65:LEU:HD13	1.96	0.48
1:A:-272:GLU:HG3	1:A:-248:LYS:O	2.15	0.47
3:L:89:GLN:HE21	3:L:96:TYR:HB3	1.79	0.47
1:A:21:THR:HA	1:A:123:LYS:O	2.15	0.46
3:L:23:CYS:SG	3:L:33:LEU:HD11	2.57	0.45
1:A:-252:ARG:HG3	1:A:-247:LEU:HD22	1.99	0.45
3:L:78:LEU:HD11	3:L:104:LEU:HD21	1.99	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:-339:ILE:HG21	1:A:-333:TYR:HB3	1.99	0.45
2:H:35:TYR:CD1	2:H:110:PRO:HG3	2.52	0.45
1:A:-4:ALA:HB2	1:A:14:ALA:HB2	1.99	0.44
1:A:84:ILE:HG13	1:A:95:ILE:HG13	1.98	0.44
1:A:-234:ILE:HA	1:A:-107:GLY:O	2.18	0.44
2:H:82:LEU:HD23	2:H:85:VAL:HG23	1.99	0.43
1:A:-331:GLY:HA3	1:A:-57:VAL:HA	2.00	0.43
1:A:-251:TYR:O	1:A:-248:LYS:HB2	2.19	0.43
3:L:19:VAL:HG21	3:L:78:LEU:HD22	2.01	0.43
2:H:161:THR:HB	2:H:209:ASN:HB3	2.01	0.43
3:L:35:TRP:CE2	3:L:73:LEU:HB2	2.54	0.42
3:L:66:GLY:HA3	3:L:71:PHE:HA	2.01	0.42
2:H:196:SER:HA	2:H:199:LEU:HD13	2.00	0.42
1:A:-252:ARG:NH1	1:A:-247:LEU:HD21	2.34	0.42
2:H:35:TYR:OH	2:H:108:ALA:O	2.28	0.42
1:A:108:TYR:N	1:A:124:VAL:O	2.53	0.41
1:A:53:GLN:NE2	1:A:55:GLY:O	2.54	0.41
1:A:-248:LYS:HB3	1:A:-246:ILE:HG23	2.03	0.41
1:A:-200:ASN:OD1	1:A:-198:GLN:N	2.54	0.41
1:A:114:HIS:ND1	1:A:114:HIS:O	2.54	0.41
2:H:4:LEU:HD23	2:H:22:CYS:SG	2.61	0.41
3:L:83:PHE:CZ	3:L:106:ILE:HG22	2.55	0.41
3:L:18:ARG:HG3	3:L:76:SER:HA	2.03	0.41
1:A:-237:LEU:HD11	1:A:-194:PHE:HA	2.03	0.40
1:A:-90:GLY:HA2	1:A:-20:MET:HE2	2.03	0.40
2:H:36:TRP:CD1	2:H:80:LEU:HB2	2.57	0.40
1:A:-253:VAL:HG21	1:A:-243:PRO:HD3	2.02	0.40
1:A:-186:ASP:O	1:A:-163:GLY:HA3	2.22	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	458/492 (93%)	440 (96%)	16 (4%)	2 (0%)	30	62
2	H	223/226 (99%)	215 (96%)	7 (3%)	1 (0%)	30	62
3	L	212/214 (99%)	206 (97%)	6 (3%)	0	100	100
All	All	893/932 (96%)	861 (96%)	29 (3%)	3 (0%)	37	67

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	H	101	TYR
1	A	106	GLY
1	A	-227	PRO

### 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	308/396 (78%)	308 (100%)	0	100	100
2	H	178/191 (93%)	178 (100%)	0	100	100
3	L	178/189 (94%)	178 (100%)	0	100	100
All	All	664/776 (86%)	664 (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. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	-97	GLN
2	H	76	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 ⓘ

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	BGC	B	1	4	12,12,12	0.54	0	17,17,17	1.09	1 (5%)
4	GLC	B	2	4	11,11,12	0.64	0	15,15,17	0.55	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	BGC	B	1	4	-	0/2/22/22	0/1/1/1
4	GLC	B	2	4	-	0/2/19/22	0/1/1/1

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	1	BGC	C3-C4-C5	2.09	114.02	110.23

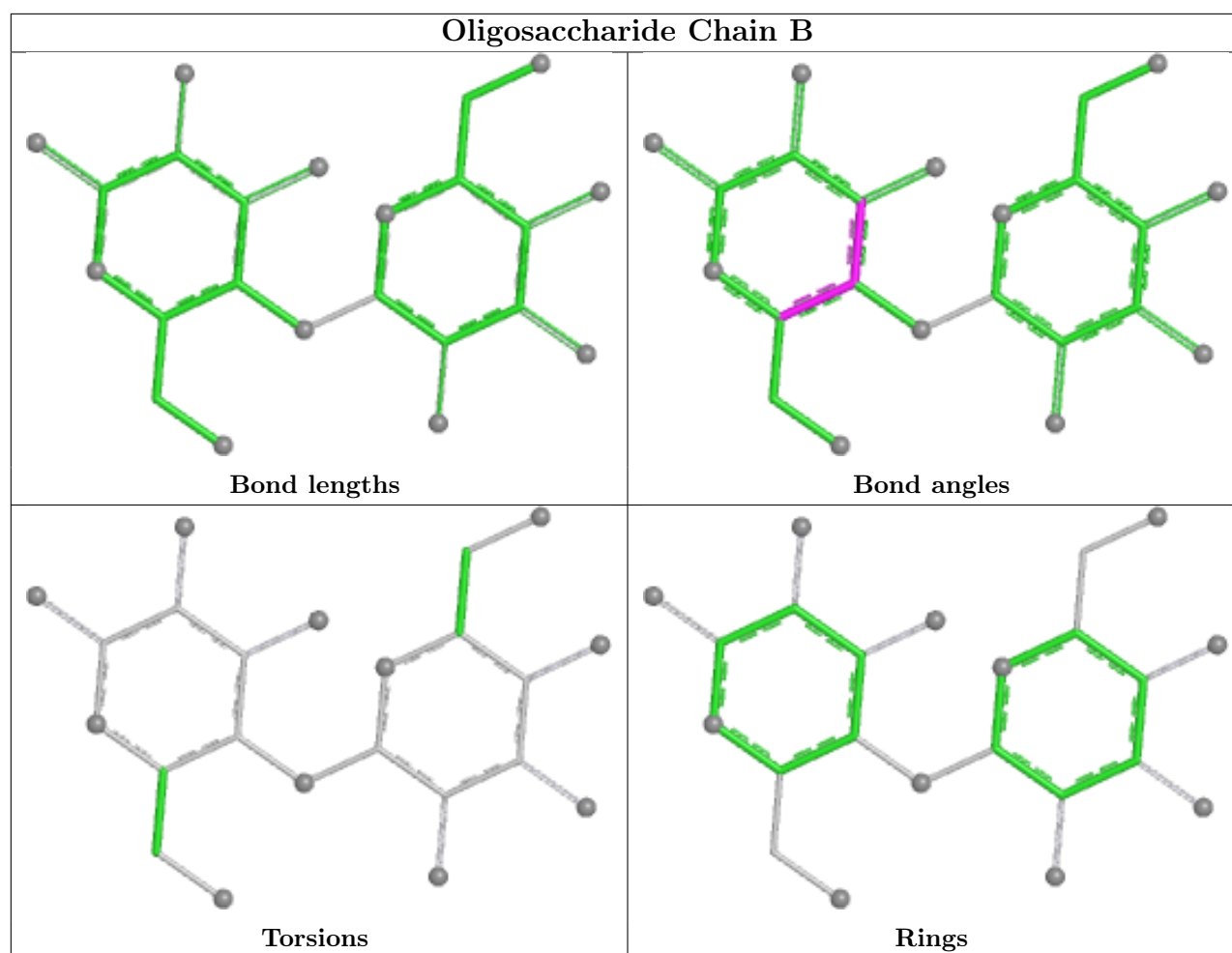
There are no chirality outliers.

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](#)

2 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	SO4	A	201	-	4,4,4	0.62	0	6,6,6	0.09	0
5	SO4	A	202	-	4,4,4	0.62	0	6,6,6	0.08	0

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 [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, 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	A	464/492 (94%)	-0.02	1 (0%) 92 84	102, 138, 178, 210	0
2	H	225/226 (99%)	-0.06	3 (1%) 74 55	78, 107, 144, 189	0
3	L	214/214 (100%)	-0.24	0 100 100	75, 105, 127, 144	0
All	All	903/932 (96%)	-0.08	4 (0%) 89 76	75, 123, 172, 210	0

All (4) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	H	140	SER	3.8
2	H	99	GLY	3.5
2	H	107	LEU	2.4
1	A	130	ALA	2.2

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

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

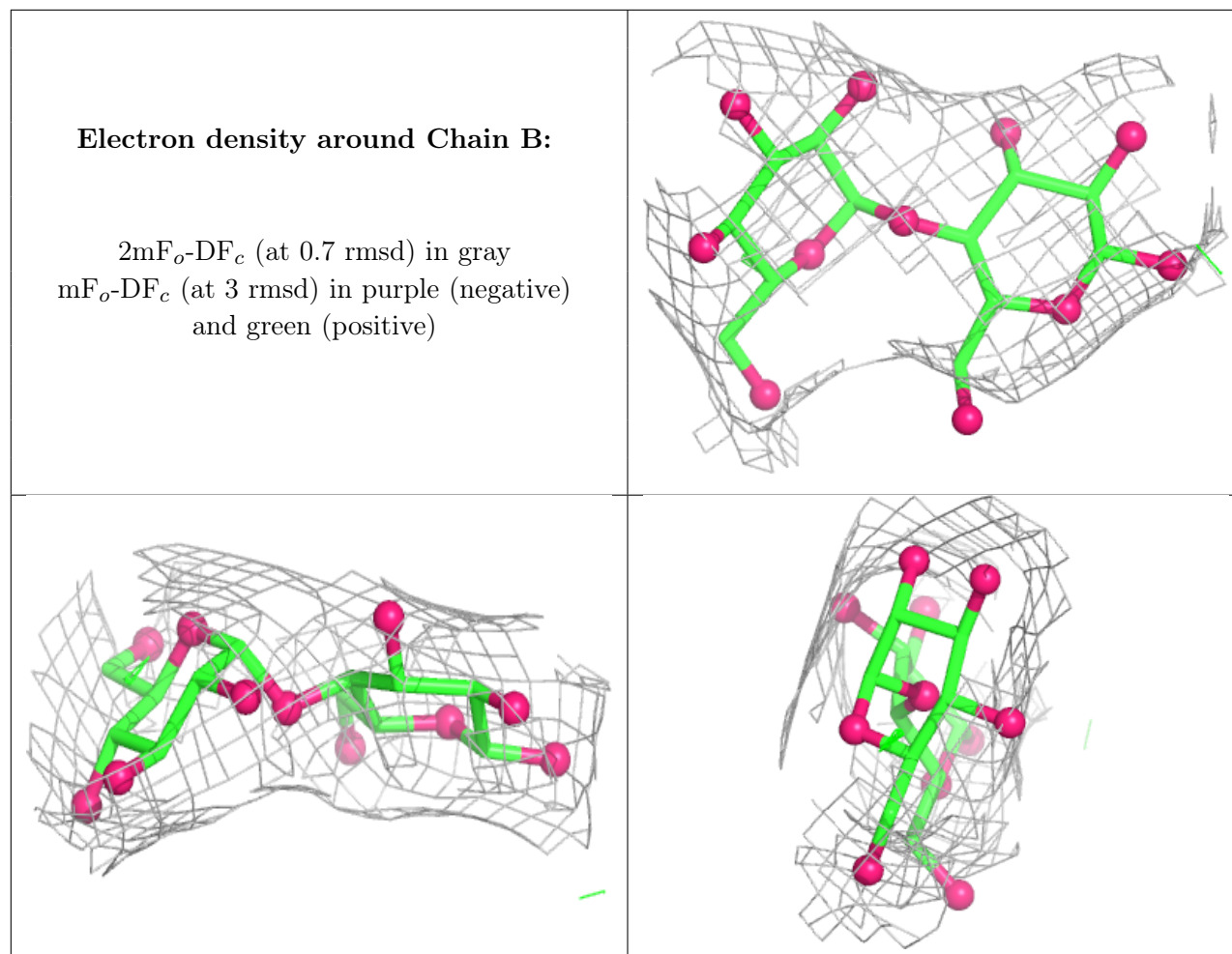
### 6.3 Carbohydrates [i](#)

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, 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
4	BGC	B	1	12/12	-	-	122,129,140,141	0
4	GLC	B	2	11/12	-	-	126,132,140,143	0

The following is a graphical depiction of the model fit to experimental electron density for oligosac-

charide. 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, 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	SO4	A	202	5/5	0.50	0.15	161,176,186,188	0
5	SO4	A	201	5/5	0.52	0.12	159,164,181,195	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.