



# wwPDB X-ray Structure Validation Summary Report ⓘ

May 22, 2020 – 02:30 pm BST

PDB ID : 3QRG  
Title : Crystal structure of antiRSVF Fab B21m  
Authors : Luo, J.; Tsui, P.; Lewandowski, F.; Spurlino, J.; Vecchio, F.D.  
Deposited on : 2011-02-17  
Resolution : 1.70 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](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.

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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 : 1.8.5 (274361), CSD as541be (2020)  
Xtriage (Phenix) : 1.13  
EDS : 2.11  
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) : 2.11

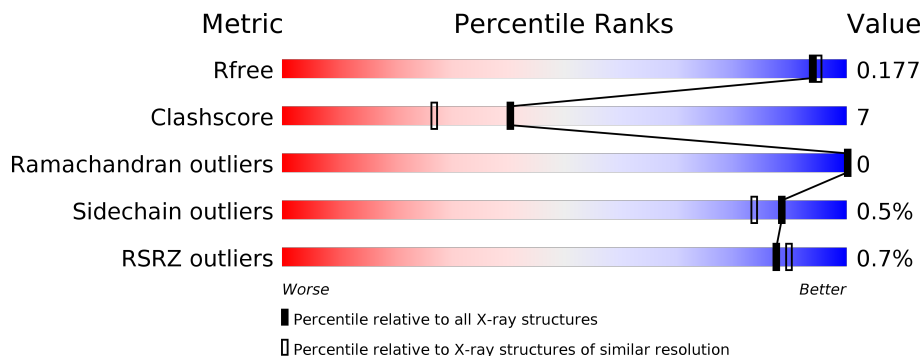
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 1.70 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))
$R_{free}$	130704	4298 (1.70-1.70)
Clashscore	141614	4695 (1.70-1.70)
Ramachandran outliers	138981	4610 (1.70-1.70)
Sidechain outliers	138945	4610 (1.70-1.70)
RSRZ outliers	127900	4222 (1.70-1.70)

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  $\geq 3$ , 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	L	218	 85% 13%
2	H	227	 89% 8%

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
4	SO4	L	230	-	-	X	-

## 2 Entry composition [i](#)

There are 6 unique types of molecules in this entry. The entry contains 4215 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.

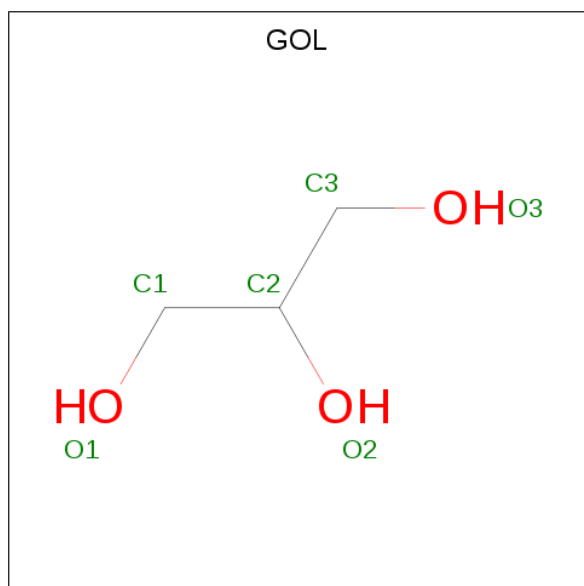
- Molecule 1 is a protein called Fab light chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	L	217	1752	1108	286	352	6	0	17	0

- Molecule 2 is a protein called Fab heavy chain Fd fragment.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	H	220	1726	1110	271	337	8	0	17	0

- Molecule 3 is GLYCEROL (three-letter code: GOL) (formula:  $C_3H_8O_3$ ).



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

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

- Molecule 4 is SULFATE ION (three-letter code: SO4) (formula: O<sub>4</sub>S).



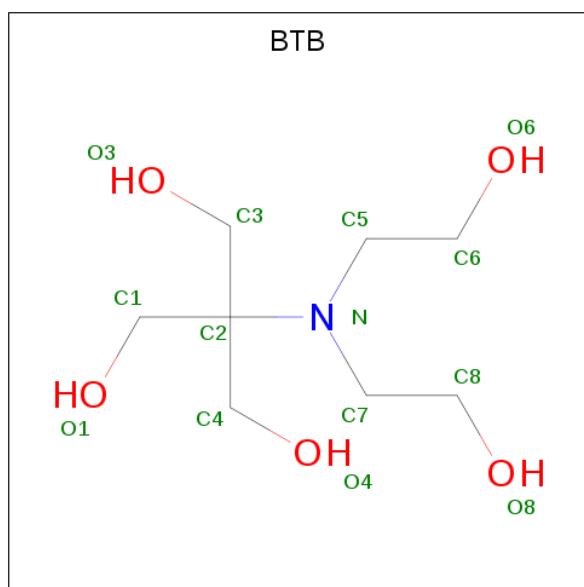
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	L	1	Total	O	S	0	0
			5	4	1		
4	L	1	Total	O	S	0	0
			5	4	1		
4	L	1	Total	O	S	0	0
			5	4	1		
4	L	1	Total	O	S	0	0
			5	4	1		

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

- Molecule 5 is 2-[BIS-(2-HYDROXY-ETHYL)-AMINO]-2-HYDROXYMETHYL-PROPAN E-1,3-DIOL (three-letter code: BTB) (formula: C<sub>8</sub>H<sub>19</sub>NO<sub>5</sub>).



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

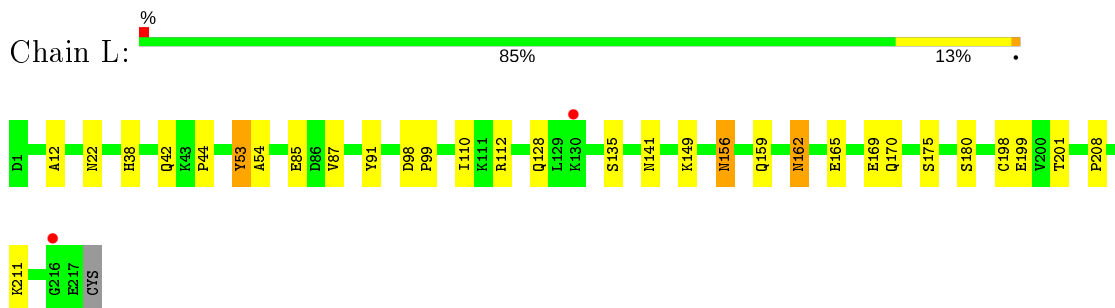
- Molecule 6 is water.

<b>Mol</b>	<b>Chain</b>	<b>Residues</b>	<b>Atoms</b>		<b>ZeroOcc</b>	<b>AltConf</b>
6	L	299	Total 299	O 299	0	0
6	H	300	Total 300	O 300	0	0

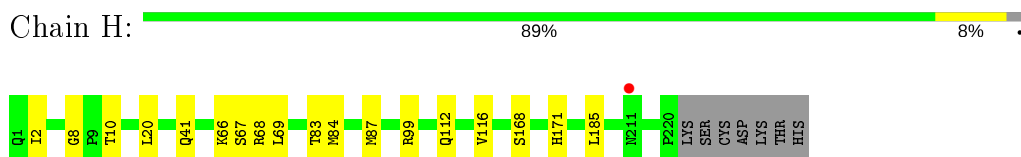
### 3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA 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: Fab light chain



- Molecule 2: Fab heavy chain Fd fragment





## 4 Data and refinement statistics

Property	Value	Source
Space group	I 2 3	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	166.41Å 166.41Å 166.41Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	41.60 – 1.70 41.60 – 1.70	Depositor EDS
% Data completeness (in resolution range)	99.1 (41.60-1.70) 99.1 (41.60-1.70)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	3.41 (at 1.70Å)	Xtrriage
Refinement program	PHENIX (phenix.refine: 1.7_650), CNX	Depositor
R, $R_{free}$	0.161 , 0.185 0.154 , 0.177	Depositor DCC
$R_{free}$ test set	4167 reflections (5.01%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	20.0	Xtrriage
Anisotropy	0.000	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.38 , 51.8	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.52$ , $\langle L^2 \rangle = 0.36$	Xtrriage
Estimated twinning fraction	0.012 for -l,-k,-h	Xtrriage
$F_o, F_c$ correlation	0.97	EDS
Total number of atoms	4215	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	23.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.22% 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 [i](#)

### 5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: GOL, PCA, SO4, BTB

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	L	1.20	6/1841 (0.3%)	1.16	2/2504 (0.1%)
2	H	1.16	1/1814 (0.1%)	1.21	3/2481 (0.1%)
All	All	1.18	7/3655 (0.2%)	1.19	5/4985 (0.1%)

The worst 5 of 7 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	L	198	CYS	CB-SG	-6.00	1.72	1.82
1	L	175	SER	CA-CB	5.55	1.61	1.52
1	L	165	GLU	CB-CG	-5.29	1.42	1.52
2	H	116	VAL	CB-CG1	5.22	1.63	1.52
1	L	53	TYR	CD1-CE1	5.10	1.47	1.39

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed( $^{\circ}$ )	Ideal( $^{\circ}$ )
1	L	112	ARG	NE-CZ-NH1	7.57	124.08	120.30
2	H	68	ARG	NE-CZ-NH1	7.36	123.98	120.30
1	L	112	ARG	NE-CZ-NH2	-6.83	116.88	120.30
2	H	68	ARG	NE-CZ-NH2	-5.32	117.64	120.30
2	H	83	THR	CA-CB-CG2	-5.01	105.38	112.40

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	L	1752	0	1748	32	0
2	H	1726	0	1763	17	0
3	H	18	0	24	5	0
3	L	36	0	47	1	0
4	H	40	0	0	0	1
4	L	30	0	0	3	0
5	H	14	0	18	3	0
6	H	300	0	0	7	1
6	L	299	0	0	6	2
All	All	4215	0	3600	52	2

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.

The worst 5 of 52 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)
2:H:99:ARG:HH22	3:H:229:GOL:H32	1.13	1.11
1:L:12:ALA:N	4:L:230:SO4:O1	1.88	1.06
1:L:42:GLN:HE22	2:H:41:GLN:HE22	1.20	0.90
2:H:67[A]:SER:OG	6:H:563:HOH:O	1.88	0.89
1:L:87[B]:VAL:HG22	1:L:110[B]:ILE:HD13	1.53	0.87

All (2) 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)
6:L:477:HOH:O	6:H:592:HOH:O[17_553]	1.91	0.29
4:H:236:SO4:O3	6:L:389:HOH:O[3_655]	2.03	0.17

## 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	L	232/218 (106%)	224 (97%)	8 (3%)	0	100	100
2	H	235/227 (104%)	230 (98%)	5 (2%)	0	100	100
All	All	467/445 (105%)	454 (97%)	13 (3%)	0	100	100

There are no Ramachandran outliers to report.

### 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	L	207/191 (108%)	205 (99%)	2 (1%)	76	67
2	H	207/197 (105%)	207 (100%)	0	100	100
All	All	414/388 (107%)	412 (100%)	2 (0%)	88	83

All (2) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	L	156	ASN
1	L	162	ASN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 12 such sidechains are listed below:

Mol	Chain	Res	Type
1	L	93	GLN
1	L	128	GLN
1	L	162	ASN
1	L	83	GLN
1	L	159	GLN

### 5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

1 non-standard protein/DNA/RNA residue is 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	PCA	H	1	2	7,8,9	1.63	1 (14%)	9,10,12	2.54	4 (44%)

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	PCA	H	1	2	-	0/0/11/13	0/1/1/1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	H	1	PCA	CD-N	4.07	1.45	1.34

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	H	1	PCA	CA-N-CD	-3.80	100.58	113.58
2	H	1	PCA	CG-CD-N	3.43	117.26	108.39
2	H	1	PCA	CB-CA-N	3.39	113.04	103.30
2	H	1	PCA	OE-CD-CG	-3.02	121.50	126.76

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

## 5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

## 5.6 Ligand geometry [i](#)

24 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
3	GOL	L	223	-	5,5,5	0.40	0	5,5,5	1.41	0
3	GOL	H	228	-	5,5,5	0.56	0	5,5,5	0.94	0
4	SO4	H	235	-	4,4,4	0.21	0	6,6,6	0.82	0
4	SO4	H	232	-	4,4,4	0.43	0	6,6,6	0.48	0
5	BTB	H	239	-	13,13,13	0.93	1 (7%)	7,16,16	2.00	3 (42%)
3	GOL	L	221	-	5,5,5	0.31	0	5,5,5	1.13	1 (20%)
4	SO4	L	228	-	4,4,4	0.23	0	6,6,6	0.55	0
4	SO4	H	234	-	4,4,4	0.42	0	6,6,6	0.71	0
4	SO4	H	237	-	4,4,4	0.26	0	6,6,6	0.64	0
4	SO4	L	226	-	4,4,4	0.29	0	6,6,6	0.54	0
4	SO4	H	238	-	4,4,4	0.23	0	6,6,6	0.36	0
3	GOL	H	229	-	5,5,5	0.45	0	5,5,5	1.45	0
4	SO4	L	230	-	4,4,4	0.42	0	6,6,6	0.70	0
3	GOL	L	222	-	5,5,5	0.41	0	5,5,5	1.06	0
3	GOL	H	230	-	5,5,5	0.50	0	5,5,5	0.92	0
4	SO4	H	231	-	4,4,4	0.81	0	6,6,6	0.46	0
3	GOL	L	220	-	5,5,5	0.56	0	5,5,5	0.83	0
3	GOL	L	224	-	5,5,5	1.35	1 (20%)	5,5,5	1.17	0
4	SO4	L	227	-	4,4,4	0.31	0	6,6,6	0.71	0
4	SO4	H	233	-	4,4,4	0.06	0	6,6,6	0.31	0
4	SO4	L	225	-	4,4,4	0.18	0	6,6,6	0.56	0
4	SO4	L	229	-	4,4,4	0.27	0	6,6,6	0.54	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
3	GOL	L	219	-	5,5,5	1.09	0	5,5,5	1.27	0
4	SO4	H	236	-	4,4,4	0.63	0	6,6,6	1.44	1 (16%)

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
3	GOL	L	223	-	-	2/4/4/4	-
3	GOL	H	228	-	-	0/4/4/4	-
3	GOL	H	230	-	-	4/4/4/4	-
3	GOL	L	220	-	-	1/4/4/4	-
3	GOL	L	224	-	-	1/4/4/4	-
3	GOL	H	229	-	-	2/4/4/4	-
5	BTB	H	239	-	-	1/21/21/21	-
3	GOL	L	221	-	-	4/4/4/4	-
3	GOL	L	222	-	-	0/4/4/4	-
3	GOL	L	219	-	-	0/4/4/4	-

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	L	224	GOL	O2-C2	-2.62	1.35	1.43
5	H	239	BTB	O1-C1	-2.15	1.35	1.42

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	H	239	BTB	O4-C4-C2	3.17	120.12	111.44
5	H	239	BTB	O1-C1-C2	-2.67	104.12	111.44
5	H	239	BTB	O3-C3-C2	2.66	118.72	111.44
4	H	236	SO4	O3-S-O1	-2.54	96.08	109.31
3	L	221	GOL	O2-C2-C3	2.01	117.96	109.12

There are no chirality outliers.

5 of 15 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	L	223	GOL	O1-C1-C2-C3

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Mol	Chain	Res	Type	Atoms
3	L	221	GOL	O1-C1-C2-C3
3	L	221	GOL	C1-C2-C3-O3
3	H	229	GOL	O1-C1-C2-C3
3	H	230	GOL	O1-C1-C2-C3

There are no ring outliers.

8 monomers are involved in 13 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	H	228	GOL	1	0
5	H	239	BTB	3	0
3	H	229	GOL	3	0
4	L	230	SO4	2	0
3	H	230	GOL	1	0
4	L	229	SO4	1	0
3	L	219	GOL	1	0
4	H	236	SO4	0	1

## 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, 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	L	217/218 (99%)	-0.27	2 (0%) 84 87	11, 20, 38, 56	0
2	H	219/227 (96%)	-0.21	1 (0%) 91 92	12, 19, 34, 48	0
All	All	436/445 (97%)	-0.24	3 (0%) 87 90	11, 19, 36, 56	0

All (3) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	L	130	LYS	3.0
2	H	211	ASN	2.6
1	L	216	GLY	2.1

### 6.2 Non-standard residues in protein, DNA, RNA chains [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
2	PCA	H	1	8/9	0.94	0.14	26,35,46,60	0

### 6.3 Carbohydrates [i](#)

There are no carbohydrates in this entry.

### 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
3	GOL	L	224	6/6	0.80	0.20	22,42,50,60	0
3	GOL	H	230	6/6	0.82	0.18	46,58,64,68	0
3	GOL	H	229	6/6	0.89	0.15	25,36,54,62	0
4	SO4	L	230	5/5	0.89	0.39	21,31,40,47	5
3	GOL	H	228	6/6	0.89	0.19	28,47,50,53	0
5	BTB	H	239	14/14	0.89	0.18	15,26,41,59	14
3	GOL	L	221	6/6	0.90	0.13	29,38,46,46	0
4	SO4	H	233	5/5	0.91	0.23	52,57,63,80	5
4	SO4	L	229	5/5	0.91	0.34	19,28,32,43	5
3	GOL	L	219	6/6	0.91	0.11	19,26,38,39	0
4	SO4	H	236	5/5	0.91	0.20	22,24,29,34	5
3	GOL	L	222	6/6	0.92	0.15	25,48,59,63	0
4	SO4	H	237	5/5	0.93	0.19	34,42,46,47	5
4	SO4	H	232	5/5	0.93	0.13	29,34,45,52	5
4	SO4	L	228	5/5	0.94	0.15	23,33,38,42	5
4	SO4	H	235	5/5	0.94	0.14	30,34,40,42	5
4	SO4	H	238	5/5	0.94	0.23	48,52,53,66	5
3	GOL	L	223	6/6	0.94	0.12	27,40,55,67	0
3	GOL	L	220	6/6	0.96	0.10	17,35,37,49	0
4	SO4	H	234	5/5	0.97	0.10	29,37,40,51	5
4	SO4	L	227	5/5	0.97	0.11	20,30,37,46	5
4	SO4	L	226	5/5	0.97	0.11	35,39,44,48	5
4	SO4	L	225	5/5	0.99	0.08	29,32,39,41	5
4	SO4	H	231	5/5	0.99	0.07	17,19,21,23	0

## 6.5 Other polymers [\(i\)](#)

There are no such residues in this entry.