



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 17, 2018 – 01:50 pm GMT

PDB ID : 1MFB  
Title : HIGH RESOLUTION STRUCTURES OF ANTIBODY FAB FRAGMENT COMPLEXED WITH CELL-SURFACE OLIGOSACCHARIDE OF PATHOGENIC SALMONELLA  
Authors : Zdanov, A.; Cygler, M.  
Deposited on : 1993-10-25  
Resolution : 2.10 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.

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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 : 1.7.3 (157068), CSD as539be (2018)  
Xtriage (Phenix) : **NOT EXECUTED**  
EDS : **NOT EXECUTED**  
Percentile statistics : 20171227.v01 (using entries in the PDB archive December 27th 2017)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : trunk30686

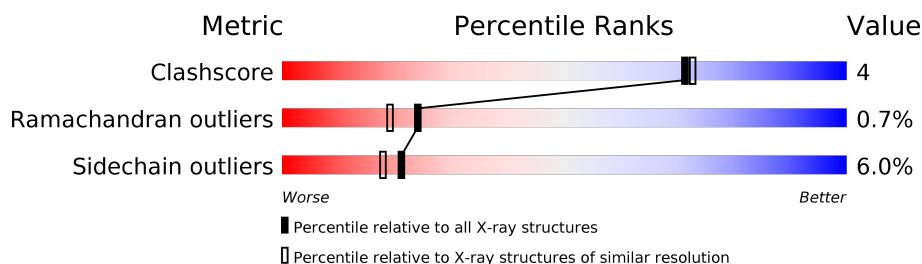
# 1 Overall quality at a glance

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))
Clashscore	122078	5107 (2.10-2.10)
Ramachandran outliers	120005	5057 (2.10-2.10)
Sidechain outliers	119972	5058 (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 on the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria. 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.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	L	215	 86% 12% ..
2	H	219	 78% 16% . .

## 2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 3505 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 IGG1-LAMBDA SE155-4 FAB (LIGHT CHAIN).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	L	212	Total	C	N	O	S	0	0	0
			1580	985	268	320	7			

There are 9 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
L	28	THR	ALA	CONFLICT	GB 387376
L	31	SER	THR	CONFLICT	GB 387376
L	32	GLY	SER	CONFLICT	GB 387376
L	34	HIS	TYR	CONFLICT	GB 387376
L	52	ASP	GLY	CONFLICT	GB 387376
L	82	PRO	THR	CONFLICT	GB 387376
L	94	CYS	TYR	CONFLICT	GB 387376
L	95	ASN	SER	CONFLICT	GB 387376
L	99	ILE	VAL	CONFLICT	GB 387376

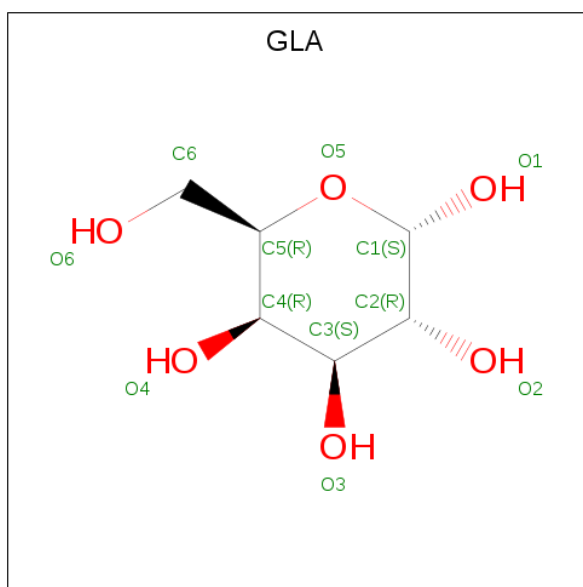
- Molecule 2 is a protein called IGG1-LAMBDA SE155-4 FAB (HEAVY CHAIN).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	H	212	Total	C	N	O	S	0	0	1
			1595	1018	266	303	8			

There are 2 discrepancies between the modelled and reference sequences:

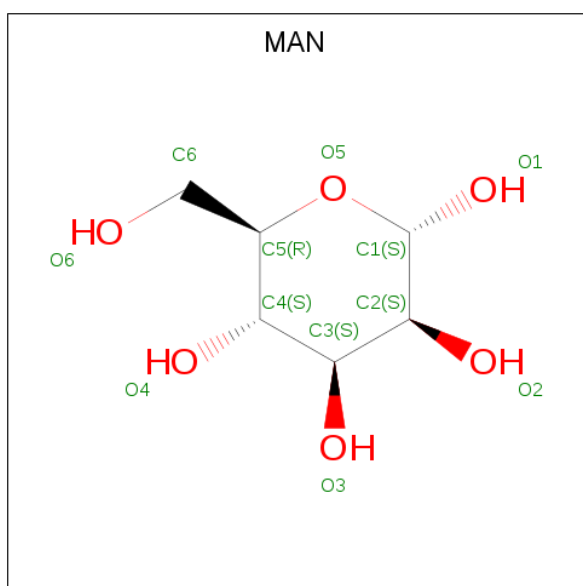
Chain	Residue	Modelled	Actual	Comment	Reference
H	?	-	ALA	DELETION	GB 208365
H	468	ARG	ASP	CONFLICT	GB 208365

- Molecule 3 is ALPHA D-GALACTOSE (three-letter code: GLA) (formula: C<sub>6</sub>H<sub>12</sub>O<sub>6</sub>).



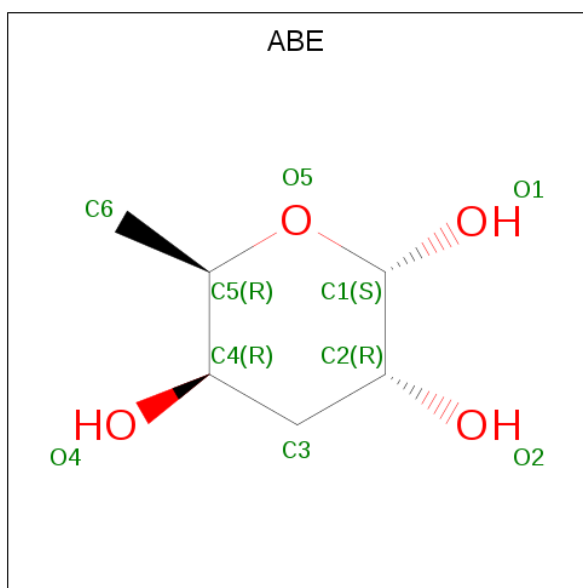
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	H	1	Total	C	O	0	0
			11	6	5		
3	H	1	Total	C	O	0	0
			11	6	5		

- Molecule 4 is ALPHA-D-MANNOSE (three-letter code: MAN) (formula:  $C_6H_{12}O_6$ ).



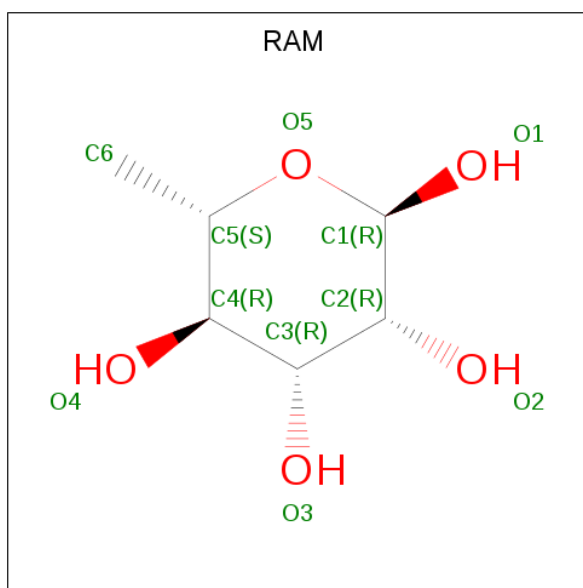
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	H	1	Total	C	O	0	0
			11	6	5		
4	H	1	Total	C	O	0	0
			11	6	5		

- Molecule 5 is ABEQUOSE (three-letter code: ABE) (formula:  $C_6H_{12}O_4$ ).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	H	1	Total	C	O	0	0
			9	6	3		

- Molecule 6 is ALPHA-L-RHAMNOSE (three-letter code: RAM) (formula:  $C_6H_{12}O_5$ ).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	H	1	Total	C	O	0	0
			11	6	5		
6	H	1	Total	C	O	0	0
			10	6	4		

- Molecule 7 is water.

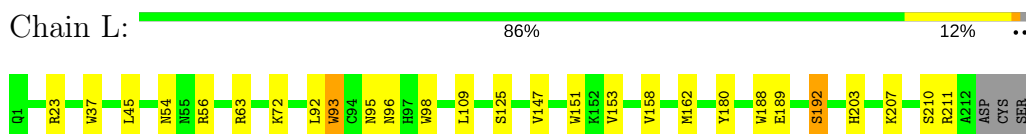
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	L	134	Total 134	O 134	0	0
7	H	122	Total 122	O 122	0	0

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

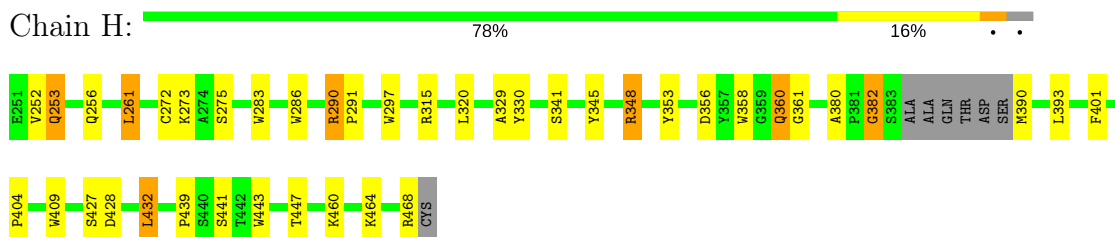
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are 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.

Note EDS was not executed.

- Molecule 1: IGG1-LAMBDA SE155-4 FAB (LIGHT CHAIN)



- Molecule 2: IGG1-LAMBDA SE155-4 FAB (HEAVY CHAIN)



## 4 Data and refinement statistics

Xtriage (Phenix) and EDS were not executed - this section is therefore incomplete.

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	47.30Å 129.00Å 79.90Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	(Not available) – 2.10	Depositor
% Data completeness (in resolution range)	(Not available) ((Not available)-2.10)	Depositor
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
Refinement program	X-PLOR	Depositor
R, $R_{free}$	0.160 , (Not available)	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	3505	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	31.0	wwPDB-VP



## 5 Model quality

### 5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: GLA, RAM, ABE, MAN

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	0.76	0/1618	1.42	17/2211 (0.8%)
2	H	0.79	0/1642	1.48	26/2246 (1.2%)
All	All	0.78	0/3260	1.45	43/4457 (1.0%)

There are no bond length outliers.

All (43) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	L	151	TRP	CD1-CG-CD2	9.08	113.56	106.30
2	H	348	ARG	NE-CZ-NH1	8.83	124.72	120.30
1	L	56	ARG	NE-CZ-NH1	8.75	124.68	120.30
2	H	297	TRP	CD1-CG-CD2	8.73	113.28	106.30
2	H	443	TRP	CD1-CG-CD2	8.67	113.24	106.30
2	H	286	TRP	CD1-CG-CD2	8.26	112.91	106.30
2	H	358	TRP	CD1-CG-CD2	8.25	112.90	106.30
2	H	283	TRP	CD1-CG-CD2	8.11	112.79	106.30
1	L	188	TRP	CD1-CG-CD2	8.08	112.76	106.30
2	H	283	TRP	CE2-CD2-CG	-8.02	100.88	107.30
1	L	98	TRP	CD1-CG-CD2	7.82	112.55	106.30
1	L	93	TRP	CD1-CG-CD2	7.78	112.53	106.30
1	L	93	TRP	CE2-CD2-CG	-7.78	101.08	107.30
2	H	358	TRP	CE2-CD2-CG	-7.75	101.10	107.30
1	L	151	TRP	CE2-CD2-CG	-7.72	101.12	107.30
2	H	348	ARG	NE-CZ-NH2	-7.71	116.45	120.30
2	H	409	TRP	CD1-CG-CD2	7.59	112.37	106.30
2	H	443	TRP	CE2-CD2-CG	-7.54	101.27	107.30
1	L	188	TRP	CE2-CD2-CG	-7.47	101.33	107.30
1	L	37	TRP	CD1-CG-CD2	7.36	112.19	106.30
2	H	286	TRP	CE2-CD2-CG	-7.25	101.50	107.30
2	H	297	TRP	CE2-CD2-CG	-7.19	101.55	107.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	L	98	TRP	CE2-CD2-CG	-6.88	101.80	107.30
1	L	37	TRP	CE2-CD2-CG	-6.70	101.94	107.30
2	H	409	TRP	CE2-CD2-CG	-6.47	102.12	107.30
1	L	93	TRP	CG-CD2-CE3	6.34	139.60	133.90
2	H	432	LEU	CA-CB-CG	6.05	129.22	115.30
2	H	358	TRP	CG-CD2-CE3	5.84	139.16	133.90
1	L	63	ARG	NE-CZ-NH1	5.79	123.20	120.30
2	H	286	TRP	CG-CD1-NE1	-5.78	104.32	110.10
1	L	93	TRP	CB-CG-CD1	-5.78	119.49	127.00
2	H	330	TYR	CB-CG-CD1	-5.76	117.55	121.00
1	L	180	TYR	CB-CG-CD1	-5.72	117.57	121.00
2	H	353	TYR	CB-CG-CD2	-5.71	117.58	121.00
2	H	297	TRP	CG-CD1-NE1	-5.52	104.58	110.10
1	L	151	TRP	CG-CD1-NE1	-5.43	104.67	110.10
2	H	443	TRP	CG-CD1-NE1	-5.38	104.72	110.10
2	H	382	GLY	N-CA-C	-5.36	99.70	113.10
2	H	315	ARG	CB-CG-CD	-5.35	97.70	111.60
2	H	315	ARG	NE-CZ-NH2	-5.29	117.66	120.30
2	H	283	TRP	CG-CD2-CE3	5.21	138.59	133.90
1	L	188	TRP	CG-CD1-NE1	-5.03	105.07	110.10
2	H	358	TRP	CB-CG-CD1	-5.02	120.48	127.00

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	L	1580	0	1511	8	0
2	H	1595	0	1547	16	0
3	H	22	0	19	0	0
4	H	22	0	17	0	0
5	H	9	0	10	0	0
6	H	21	0	20	0	0
7	H	122	0	0	1	0
7	L	134	0	0	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
All	All	3505	0	3124	24	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 4.

All (24) 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:253:GLN:HB2	2:H:275:SER:HB2	1.75	0.69
2:H:256:GLN:HE22	2:H:345:TYR:HA	1.62	0.63
2:H:256:GLN:HG2	2:H:272:CYS:SG	2.41	0.61
1:L:23:ARG:HH11	1:L:72:LYS:HD3	1.67	0.59
1:L:23:ARG:HB3	1:L:72:LYS:HG2	1.87	0.57
1:L:192:SER:O	1:L:210:SER:HA	2.09	0.53
2:H:390:MET:HG2	2:H:439:PRO:HA	1.91	0.51
2:H:380:ALA:HB3	2:H:468:ARG:HG3	1.94	0.49
2:H:256:GLN:NE2	2:H:361:GLY:H	2.11	0.48
2:H:360:GLN:NE2	2:H:360:GLN:H	2.12	0.48
1:L:95:ASN:HB2	7:L:830:HOH:O	2.13	0.48
2:H:447:THR:HG23	2:H:464:LYS:HG3	1.97	0.47
2:H:290:ARG:HH21	2:H:341:SER:HB2	1.78	0.47
2:H:252:VAL:HG21	2:H:348:ARG:NH2	2.33	0.44
1:L:153:VAL:HG23	1:L:158:VAL:HG21	2.00	0.44
2:H:329:ALA:HB1	7:H:657:HOH:O	2.17	0.43
1:L:23:ARG:NH1	1:L:72:LYS:HD3	2.32	0.43
1:L:93:TRP:CZ2	1:L:96:ASN:HA	2.54	0.43
2:H:252:VAL:HA	2:H:275:SER:O	2.19	0.42
2:H:360:GLN:H	2:H:360:GLN:HE21	1.67	0.42
2:H:382:GLY:HA2	2:H:468:ARG:HB2	2.00	0.42
2:H:261:LEU:HD21	2:H:401:PHE:HZ	1.85	0.42
2:H:348:ARG:HD2	2:H:356:ASP:OD1	2.20	0.41
1:L:189:GLU:HA	1:L:211:ARG:HD2	2.03	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	L	210/215 (98%)	202 (96%)	7 (3%)	1 (0%)	31	27
2	H	208/219 (95%)	196 (94%)	10 (5%)	2 (1%)	17	12
All	All	418/434 (96%)	398 (95%)	17 (4%)	3 (1%)	24	19

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	L	109	LEU
2	H	291	PRO
2	H	427	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	L	176/183 (96%)	167 (95%)	9 (5%)	26	24
2	H	175/184 (95%)	163 (93%)	12 (7%)	17	13
All	All	351/367 (96%)	330 (94%)	21 (6%)	21	18

All (21) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	L	45	LEU
1	L	54	ASN
1	L	92	LEU

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Mol	Chain	Res	Type
1	L	125	SER
1	L	147	VAL
1	L	162	MET
1	L	192	SER
1	L	203	HIS
1	L	207	LYS
2	H	253	GLN
2	H	261	LEU
2	H	273	LYS
2	H	290	ARG
2	H	320	LEU
2	H	360	GLN
2	H	393	LEU
2	H	404	PRO
2	H	428	ASP
2	H	432	LEU
2	H	441	SER
2	H	460	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	L	36	ASN
1	L	54	ASN
1	L	191	HIS
1	L	197	GLN
2	H	256	GLN
2	H	289	GLN
2	H	305	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 [i](#)

There are no carbohydrates in this entry.

## 5.6 Ligand geometry [i](#)

7 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	GLA	H	511	4,6	11,11,12	0.69	0	15,15,17	1.18	2 (13%)
4	MAN	H	512	3,5,6	11,11,12	0.64	0	15,15,17	1.64	1 (6%)
5	ABE	H	513	4	8,9,10	1.81	2 (25%)	9,12,14	0.87	0
6	RAM	H	514	4	11,11,11	1.01	0	15,16,16	1.10	1 (6%)
3	GLA	H	521	4	11,11,12	1.30	1 (9%)	15,15,17	1.18	2 (13%)
4	MAN	H	522	3,6	11,11,12	0.73	0	15,15,17	1.43	3 (20%)
6	RAM	H	524	3,4	9,10,11	0.79	0	13,14,16	1.18	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
3	GLA	H	511	4,6	-	0/2/19/22	0/1/1/1
4	MAN	H	512	3,5,6	-	0/2/19/22	0/1/1/1
5	ABE	H	513	4	-	0/0/13/16	0/1/1/1
6	RAM	H	514	4	-	0/0/20/20	0/1/1/1
3	GLA	H	521	4	-	0/2/19/22	0/1/1/1
4	MAN	H	522	3,6	-	0/2/19/22	0/1/1/1
6	RAM	H	524	3,4	-	0/0/17/20	0/1/1/1

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	H	513	ABE	C4-C5	3.00	1.58	1.52
3	H	521	GLA	C2-C3	3.16	1.57	1.52
5	H	513	ABE	C1-C2	3.77	1.54	1.51

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	H	522	MAN	O2-C2-C1	-2.24	104.65	109.17
3	H	521	GLA	O4-C4-C3	-2.07	105.52	110.34
3	H	511	GLA	C6-C5-C4	2.06	117.85	112.99
4	H	522	MAN	C1-C2-C3	2.13	112.36	109.66
3	H	521	GLA	C6-C5-C4	2.30	118.42	112.99
6	H	514	RAM	C6-C5-C4	2.69	117.69	113.06
3	H	511	GLA	C1-O5-C5	3.07	116.41	112.19
4	H	522	MAN	C1-O5-C5	4.10	117.82	112.19
4	H	512	MAN	C1-O5-C5	5.54	119.81	112.19

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 [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

EDS was not executed - this section is therefore empty.

### 6.2 Non-standard residues in protein, DNA, RNA chains

EDS was not executed - this section is therefore empty.

### 6.3 Carbohydrates

EDS was not executed - this section is therefore empty.

### 6.4 Ligands

EDS was not executed - this section is therefore empty.

### 6.5 Other polymers

EDS was not executed - this section is therefore empty.