



# wwPDB X-ray Structure Validation Summary Report ⓘ

May 13, 2020 – 07:27 am BST

PDB ID : 5MVZ  
Title : Fab 4AB007 bound to Interleukin-1-beta  
Authors : Stark, W.; Seibert, V.  
Deposited on : 2017-01-17  
Resolution : 2.15 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

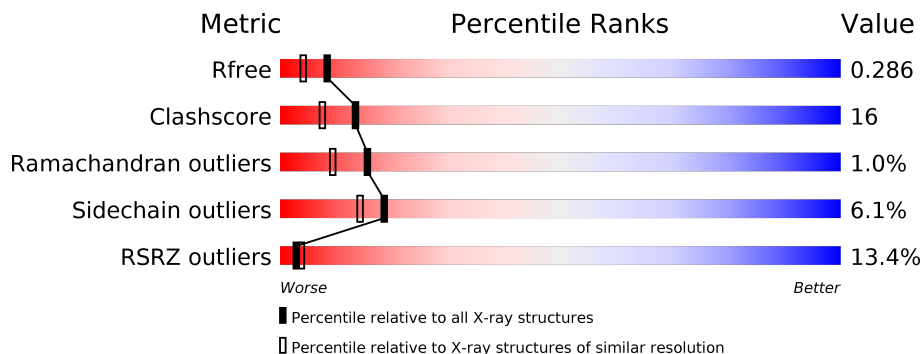
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.15 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	1479 (2.16-2.16)
Clashscore	141614	1585 (2.16-2.16)
Ramachandran outliers	138981	1560 (2.16-2.16)
Sidechain outliers	138945	1559 (2.16-2.16)
RSRZ outliers	127900	1456 (2.16-2.16)

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	A	240	
1	H	240	
2	B	249	
2	L	249	
3	U	177	
3	V	177	

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:

<b>Mol</b>	<b>Type</b>	<b>Chain</b>	<b>Res</b>	<b>Chirality</b>	<b>Geometry</b>	<b>Clashes</b>	<b>Electron density</b>
4	GOL	A	301	-	-	X	-
4	GOL	H	301	-	-	-	X
4	GOL	L	301	-	-	X	-

## 2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 9536 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 4AB007 H-chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	H	221	Total 1659	C 1053	N 269	O 328	S 9	0	0	0
1	A	221	Total 1659	C 1053	N 269	O 328	S 9	0	0	0

- Molecule 2 is a protein called Fab 4AB007 L-chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	L	215	Total 1662	C 1040	N 282	O 335	S 5	0	0	0
2	B	214	Total 1653	C 1035	N 281	O 332	S 5	0	0	0

- Molecule 3 is a protein called Interleukin-1 beta.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
3	U	136	Total 1084	C 694	N 175	O 208	S 7	0	0	0
3	V	144	Total 1137	C 727	N 187	O 215	S 8	0	0	0

There are 48 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
U	-23	MET	-	initiating methionine	UNP B5BUQ8
U	-22	ALA	-	expression tag	UNP B5BUQ8
U	-21	HIS	-	expression tag	UNP B5BUQ8
U	-20	HIS	-	expression tag	UNP B5BUQ8
U	-19	HIS	-	expression tag	UNP B5BUQ8
U	-18	HIS	-	expression tag	UNP B5BUQ8
U	-17	HIS	-	expression tag	UNP B5BUQ8
U	-16	HIS	-	expression tag	UNP B5BUQ8

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Chain	Residue	Modelled	Actual	Comment	Reference
U	-15	HIS	-	expression tag	UNP B5BUQ8
U	-14	HIS	-	expression tag	UNP B5BUQ8
U	-13	HIS	-	expression tag	UNP B5BUQ8
U	-12	HIS	-	expression tag	UNP B5BUQ8
U	-11	HIS	-	expression tag	UNP B5BUQ8
U	-10	HIS	-	expression tag	UNP B5BUQ8
U	-9	PRO	-	expression tag	UNP B5BUQ8
U	-8	ALA	-	expression tag	UNP B5BUQ8
U	-7	GLY	-	expression tag	UNP B5BUQ8
U	-6	GLU	-	expression tag	UNP B5BUQ8
U	-5	ASN	-	expression tag	UNP B5BUQ8
U	-4	LEU	-	expression tag	UNP B5BUQ8
U	-3	TYR	-	expression tag	UNP B5BUQ8
U	-2	PHE	-	expression tag	UNP B5BUQ8
U	-1	GLN	-	expression tag	UNP B5BUQ8
U	0	GLY	-	expression tag	UNP B5BUQ8
V	-23	MET	-	initiating methionine	UNP B5BUQ8
V	-22	ALA	-	expression tag	UNP B5BUQ8
V	-21	HIS	-	expression tag	UNP B5BUQ8
V	-20	HIS	-	expression tag	UNP B5BUQ8
V	-19	HIS	-	expression tag	UNP B5BUQ8
V	-18	HIS	-	expression tag	UNP B5BUQ8
V	-17	HIS	-	expression tag	UNP B5BUQ8
V	-16	HIS	-	expression tag	UNP B5BUQ8
V	-15	HIS	-	expression tag	UNP B5BUQ8
V	-14	HIS	-	expression tag	UNP B5BUQ8
V	-13	HIS	-	expression tag	UNP B5BUQ8
V	-12	HIS	-	expression tag	UNP B5BUQ8
V	-11	HIS	-	expression tag	UNP B5BUQ8
V	-10	HIS	-	expression tag	UNP B5BUQ8
V	-9	PRO	-	expression tag	UNP B5BUQ8
V	-8	ALA	-	expression tag	UNP B5BUQ8
V	-7	GLY	-	expression tag	UNP B5BUQ8
V	-6	GLU	-	expression tag	UNP B5BUQ8
V	-5	ASN	-	expression tag	UNP B5BUQ8
V	-4	LEU	-	expression tag	UNP B5BUQ8
V	-3	TYR	-	expression tag	UNP B5BUQ8
V	-2	PHE	-	expression tag	UNP B5BUQ8
V	-1	GLN	-	expression tag	UNP B5BUQ8
V	0	GLY	-	expression tag	UNP B5BUQ8

- Molecule 4 is GLYCEROL (three-letter code: GOL) (formula: C<sub>3</sub>H<sub>8</sub>O<sub>3</sub>).



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

- Molecule 5 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	H	136	Total O 136 136	0	0
5	L	152	Total O 152 152	0	0
5	U	28	Total O 28 28	0	0
5	A	130	Total O 130 130	0	0

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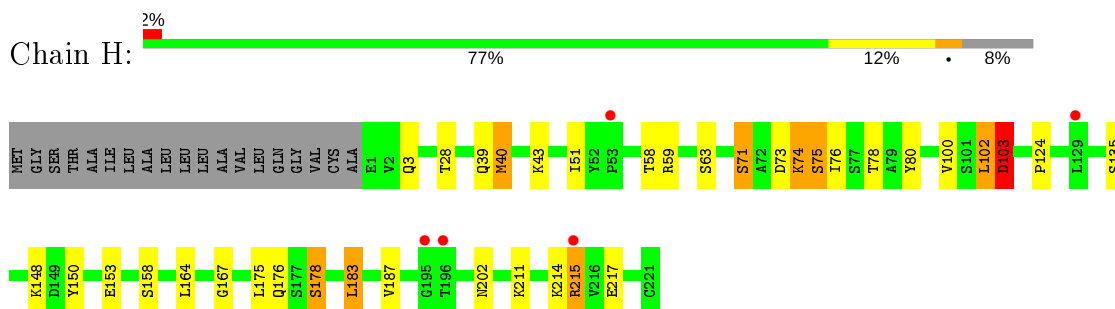
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<b>Mol</b>	<b>Chain</b>	<b>Residues</b>	<b>Atoms</b>		<b>ZeroOcc</b>	<b>AltConf</b>
5	B	156	Total 156	O 156	0	0
5	V	32	Total 32	O 32	0	0

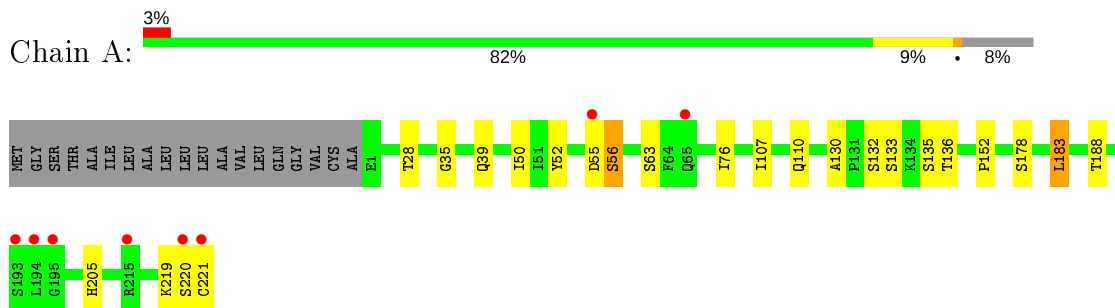
### 3 Residue-property plots [i](#)

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.

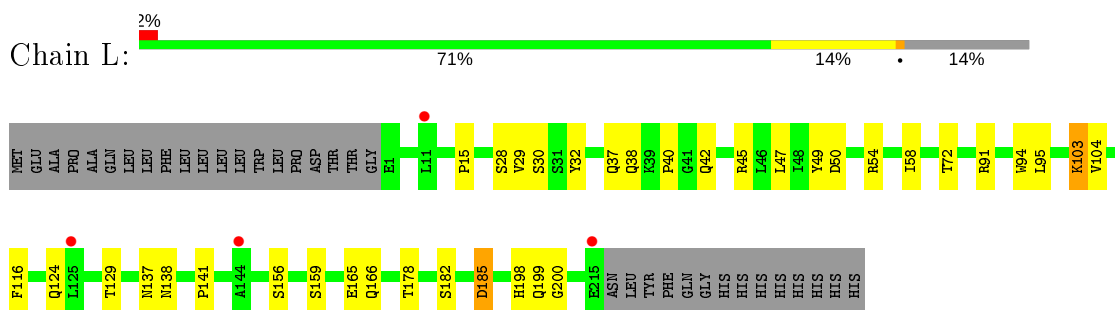
- Molecule 1: Fab 4AB007 H-chain



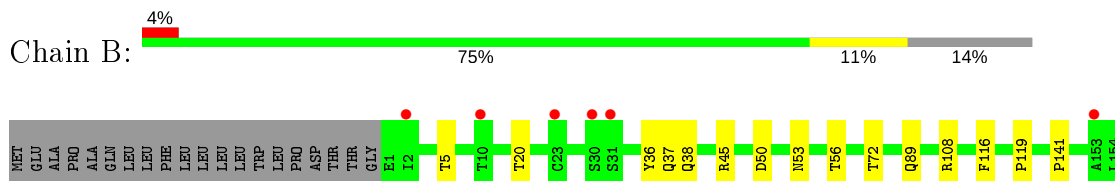
- Molecule 1: Fab 4AB007 H-chain



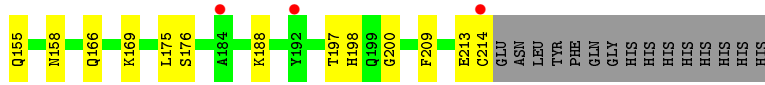
- Molecule 2: Fab 4AB007 L-chain



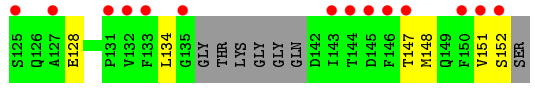
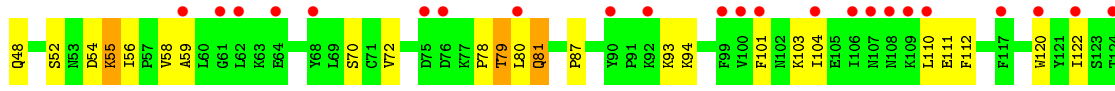
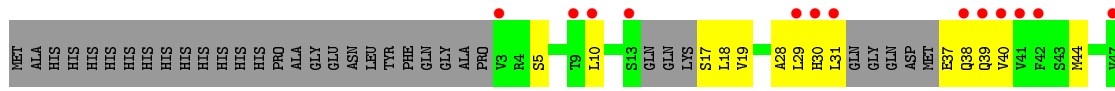
- Molecule 2: Fab 4AB007 L-chain



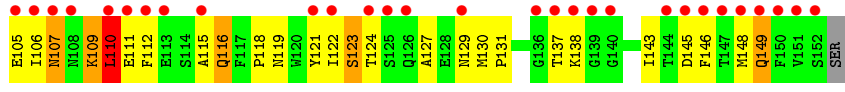
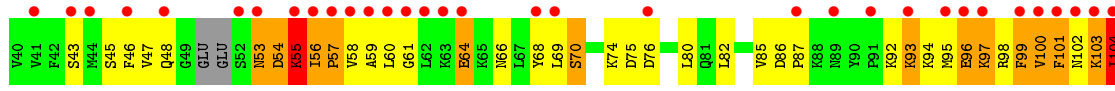
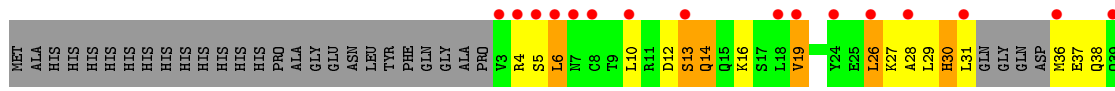




● Molecule 3: Interleukin-1 beta



● Molecule 3: Interleukin-1 beta



## 4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	111.04Å 158.38Å 153.44Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	47.68 – 2.15 45.53 – 2.15	Depositor EDS
% Data completeness (in resolution range)	99.6 (47.68-2.15) 99.6 (45.53-2.15)	Depositor EDS
$R_{merge}$	0.07	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.09 (at 2.16Å)	Xtrriage
Refinement program	REFMAC 5.8.0071	Depositor
R, $R_{free}$	0.219 , 0.281 0.232 , 0.286	Depositor DCC
$R_{free}$ test set	3665 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	40.1	Xtrriage
Anisotropy	0.025	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.35 , 73.2	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.45$ , $\langle L^2 \rangle = 0.27$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
$F_o, F_c$ correlation	0.93	EDS
Total number of atoms	9536	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	66.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 27.28 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 2.2605e-03.*

<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

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	0.71	0/1702	0.76	1/2316 (0.0%)
1	H	0.71	0/1702	0.75	1/2316 (0.0%)
2	B	0.75	0/1690	0.85	1/2297 (0.0%)
2	L	0.77	0/1699	0.84	0/2309
3	U	0.48	0/1103	0.65	0/1484
3	V	0.64	0/1157	0.90	2/1556 (0.1%)
All	All	0.70	0/9053	0.80	5/12278 (0.0%)

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
3	V	0	2

There are no bond length outliers.

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	V	104	ILE	CG1-CB-CG2	8.56	130.24	111.40
3	V	110	LEU	CA-CB-CG	6.87	131.10	115.30
2	B	108	ARG	NE-CZ-NH2	-5.74	117.43	120.30
1	H	103	ASP	CB-CG-OD2	5.63	123.37	118.30
1	A	183	LEU	CA-CB-CG	5.55	128.06	115.30

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
3	V	137	THR	Peptide
3	V	57	PRO	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	1659	0	1626	21	9
1	H	1659	0	1625	28	0
2	B	1653	0	1603	17	0
2	L	1662	0	1608	36	0
3	U	1084	0	1066	38	0
3	V	1137	0	1124	155	10
4	A	6	0	8	4	0
4	B	18	0	24	1	0
4	H	6	0	8	3	0
4	L	18	0	24	9	0
5	A	130	0	0	0	0
5	B	156	0	0	3	1
5	H	136	0	0	7	0
5	L	152	0	0	10	1
5	U	28	0	0	2	1
5	V	32	0	0	11	0
All	All	9536	0	8716	284	12

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 16.

The worst 5 of 284 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)
3:V:4:ARG:CZ	5:V:201:HOH:O	1.75	1.26
3:V:45:SER:O	3:V:59:ALA:HB3	1.31	1.24
3:V:104:ILE:HD12	3:V:110:LEU:CA	1.72	1.17
3:V:104:ILE:CD1	3:V:109:LYS:O	1.93	1.16
3:V:47:VAL:HG22	3:V:58:VAL:CG1	1.77	1.14

The worst 5 of 12 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)
1:A:132:SER:O	3:V:55:LYS:NZ[5_555]	1.12	1.08
1:A:133:SER:CA	3:V:55:LYS:CE[5_555]	1.33	0.87
1:A:133:SER:N	3:V:55:LYS:CE[5_555]	1.61	0.59
1:A:133:SER:CA	3:V:55:LYS:CD[5_555]	1.67	0.53
1:A:136:THR:OG1	3:V:55:LYS:NZ[5_555]	1.90	0.30

## 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	219/240 (91%)	214 (98%)	3 (1%)	2 (1%)	17	11
1	H	219/240 (91%)	209 (95%)	8 (4%)	2 (1%)	17	11
2	B	212/249 (85%)	203 (96%)	9 (4%)	0	100	100
2	L	213/249 (86%)	204 (96%)	9 (4%)	0	100	100
3	U	128/177 (72%)	123 (96%)	5 (4%)	0	100	100
3	V	138/177 (78%)	114 (83%)	17 (12%)	7 (5%)	2	0
All	All	1129/1332 (85%)	1067 (94%)	51 (4%)	11 (1%)	15	9

5 of 11 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
3	V	103	LYS
1	H	178	SER
3	V	101	PHE
3	V	138	LYS
1	A	56	SER

### 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	188/201 (94%)	185 (98%)	3 (2%)	62	67
1	H	188/201 (94%)	176 (94%)	12 (6%)	17	12
2	B	186/217 (86%)	182 (98%)	4 (2%)	52	55
2	L	187/217 (86%)	183 (98%)	4 (2%)	53	57
3	U	125/160 (78%)	115 (92%)	10 (8%)	12	7
3	V	128/160 (80%)	100 (78%)	28 (22%)	1	0
All	All	1002/1156 (87%)	941 (94%)	61 (6%)	18	14

5 of 61 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	107	ILE
3	V	13	SER
3	V	110	LEU
1	A	183	LEU
2	B	169	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 22 such sidechains are listed below:

Mol	Chain	Res	Type
1	A	169	HIS
2	B	37	GLN
3	V	107	ASN
1	A	176	GLN
1	A	205	HIS

### 5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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](#)

8 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	GOL	B	303	-	5,5,5	0.24	0	5,5,5	0.57	0
4	GOL	L	303	-	5,5,5	0.44	0	5,5,5	0.27	0
4	GOL	A	301	-	5,5,5	0.71	0	5,5,5	1.44	1 (20%)
4	GOL	B	301	-	5,5,5	0.67	0	5,5,5	0.89	0
4	GOL	L	301	-	5,5,5	0.60	0	5,5,5	1.22	1 (20%)
4	GOL	B	302	-	5,5,5	0.67	0	5,5,5	0.67	0
4	GOL	L	302	-	5,5,5	0.61	0	5,5,5	0.97	0
4	GOL	H	301	-	5,5,5	0.27	0	5,5,5	0.41	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	GOL	B	303	-	-	4/4/4/4	-
4	GOL	L	303	-	-	2/4/4/4	-
4	GOL	A	301	-	-	0/4/4/4	-
4	GOL	B	301	-	-	0/4/4/4	-
4	GOL	L	301	-	-	2/4/4/4	-
4	GOL	B	302	-	-	2/4/4/4	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	GOL	L	302	-	-	2/4/4/4	-
4	GOL	H	301	-	-	2/4/4/4	-

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	301	GOL	O2-C2-C1	-2.09	99.92	109.12
4	L	301	GOL	O3-C3-C2	2.05	120.03	110.20

There are no chirality outliers.

5 of 14 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	L	303	GOL	O1-C1-C2-C3
4	L	301	GOL	O1-C1-C2-C3
4	B	302	GOL	O1-C1-C2-C3
4	H	301	GOL	C1-C2-C3-O3
4	B	302	GOL	O1-C1-C2-O2

There are no ring outliers.

5 monomers are involved in 17 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	301	GOL	4	0
4	L	301	GOL	7	0
4	B	302	GOL	1	0
4	L	302	GOL	2	0
4	H	301	GOL	3	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, 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	221/240 (92%)	0.43	8 (3%) 42 51	45, 59, 77, 94	0
1	H	221/240 (92%)	0.32	5 (2%) 60 68	48, 61, 79, 91	0
2	B	214/249 (85%)	0.59	9 (4%) 36 45	40, 55, 79, 108	0
2	L	215/249 (86%)	0.54	4 (1%) 66 74	41, 54, 76, 125	0
3	U	136/177 (76%)	1.74	50 (36%) 0 0	51, 87, 107, 125	0
3	V	144/177 (81%)	2.89	78 (54%) 0 0	60, 86, 116, 131	0
All	All	1151/1332 (86%)	0.92	154 (13%) 3 4	40, 62, 100, 131	0

The worst 5 of 154 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
3	V	110	LEU	11.5
3	V	100	VAL	11.0
3	V	3	VAL	9.4
3	V	58	VAL	9.4
3	V	18	LEU	8.7

### 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](#)

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
4	GOL	H	301	6/6	0.67	0.55	70,75,82,87	0
4	GOL	L	301	6/6	0.75	0.28	49,56,68,69	0
4	GOL	L	303	6/6	0.82	0.15	60,73,76,79	0
4	GOL	A	301	6/6	0.82	0.23	46,59,63,70	0
4	GOL	L	302	6/6	0.84	0.21	64,65,66,78	0
4	GOL	B	302	6/6	0.85	0.25	48,57,59,66	0
4	GOL	B	303	6/6	0.90	0.21	65,74,78,82	0
4	GOL	B	301	6/6	0.94	0.15	53,54,59,62	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.