



# wwPDB X-ray Structure Validation Summary Report ⓘ

Feb 19, 2016 – 09:53 PM GMT

PDB ID : 5BOB  
Title : Crystal Structure of the Meningitis Pathogen Streptococcus suis adhesion Fhb  
Authors : Jiang, Y.; Zhang, C.; Yu, Y.  
Deposited on : 2015-05-27  
Resolution : 1.50 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  
<http://wwpdb.org/validation/2016/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 : 1.7.1 (RC1), CSD as537be (2016)  
Xtriage (Phenix) : 1.9-1692  
EDS : rb-20026982  
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)  
Refmac : 5.8.0135  
CCP4 : 6.5.0  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : rb-20026982

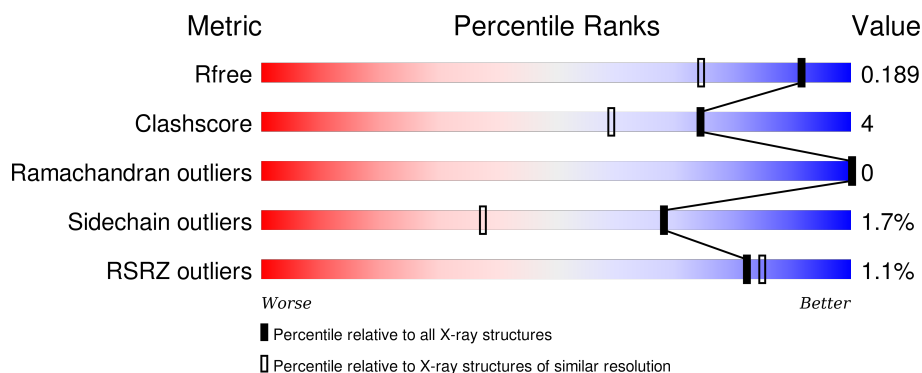
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 1.50 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))
$R_{free}$	91344	2072 (1.50-1.50)
Clashscore	102246	2274 (1.50-1.50)
Ramachandran outliers	100387	2218 (1.50-1.50)
Sidechain outliers	100360	2216 (1.50-1.50)
RSRZ outliers	91569	2075 (1.50-1.50)

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.

Mol	Chain	Length	Quality of chain
1	A	226	
1	B	226	
1	C	226	
1	D	226	
1	E	226	

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
2	GOL	A	402	-	-	-	X
2	GOL	B	401	-	-	-	X
2	GOL	C	401	-	-	-	X
2	GOL	D	401	-	-	-	X
2	GOL	D	402	-	-	-	X
2	GOL	E	402	-	-	-	X
2	GOL	E	403	-	-	-	X

## 2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 9071 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 Translation initiation factor 2 (IF-2 GTPase).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	203	Total	C	N	O	Se	0	1	0
			1606	1008	272	325	1			
1	B	199	Total	C	N	O	Se	0	0	0
			1579	993	268	317	1			
1	C	198	Total	C	N	O	Se	0	1	0
			1579	995	267	316	1			
1	D	200	Total	C	N	O	Se	0	1	0
			1590	999	270	320	1			
1	E	196	Total	C	N	O	Se	0	1	0
			1573	989	267	316	1			

There are 105 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	118	MSE	-	expression tag	UNP A4VT01
A	119	GLY	-	expression tag	UNP A4VT01
A	120	SER	-	expression tag	UNP A4VT01
A	121	SER	-	expression tag	UNP A4VT01
A	122	HIS	-	expression tag	UNP A4VT01
A	123	HIS	-	expression tag	UNP A4VT01
A	124	HIS	-	expression tag	UNP A4VT01
A	125	HIS	-	expression tag	UNP A4VT01
A	126	HIS	-	expression tag	UNP A4VT01
A	127	HIS	-	expression tag	UNP A4VT01
A	128	SER	-	expression tag	UNP A4VT01
A	129	SER	-	expression tag	UNP A4VT01
A	130	GLY	-	expression tag	UNP A4VT01
A	131	LEU	-	expression tag	UNP A4VT01
A	132	VAL	-	expression tag	UNP A4VT01
A	133	PRO	-	expression tag	UNP A4VT01
A	134	ARG	-	expression tag	UNP A4VT01
A	135	GLY	-	expression tag	UNP A4VT01
A	136	SER	-	expression tag	UNP A4VT01

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Chain	Residue	Modelled	Actual	Comment	Reference
A	137	HIS	-	expression tag	UNP A4VT01
A	138	MSE	-	expression tag	UNP A4VT01
B	118	MSE	-	expression tag	UNP A4VT01
B	119	GLY	-	expression tag	UNP A4VT01
B	120	SER	-	expression tag	UNP A4VT01
B	121	SER	-	expression tag	UNP A4VT01
B	122	HIS	-	expression tag	UNP A4VT01
B	123	HIS	-	expression tag	UNP A4VT01
B	124	HIS	-	expression tag	UNP A4VT01
B	125	HIS	-	expression tag	UNP A4VT01
B	126	HIS	-	expression tag	UNP A4VT01
B	127	HIS	-	expression tag	UNP A4VT01
B	128	SER	-	expression tag	UNP A4VT01
B	129	SER	-	expression tag	UNP A4VT01
B	130	GLY	-	expression tag	UNP A4VT01
B	131	LEU	-	expression tag	UNP A4VT01
B	132	VAL	-	expression tag	UNP A4VT01
B	133	PRO	-	expression tag	UNP A4VT01
B	134	ARG	-	expression tag	UNP A4VT01
B	135	GLY	-	expression tag	UNP A4VT01
B	136	SER	-	expression tag	UNP A4VT01
B	137	HIS	-	expression tag	UNP A4VT01
B	138	MSE	-	expression tag	UNP A4VT01
C	118	MSE	-	expression tag	UNP A4VT01
C	119	GLY	-	expression tag	UNP A4VT01
C	120	SER	-	expression tag	UNP A4VT01
C	121	SER	-	expression tag	UNP A4VT01
C	122	HIS	-	expression tag	UNP A4VT01
C	123	HIS	-	expression tag	UNP A4VT01
C	124	HIS	-	expression tag	UNP A4VT01
C	125	HIS	-	expression tag	UNP A4VT01
C	126	HIS	-	expression tag	UNP A4VT01
C	127	HIS	-	expression tag	UNP A4VT01
C	128	SER	-	expression tag	UNP A4VT01
C	129	SER	-	expression tag	UNP A4VT01
C	130	GLY	-	expression tag	UNP A4VT01
C	131	LEU	-	expression tag	UNP A4VT01
C	132	VAL	-	expression tag	UNP A4VT01
C	133	PRO	-	expression tag	UNP A4VT01
C	134	ARG	-	expression tag	UNP A4VT01
C	135	GLY	-	expression tag	UNP A4VT01
C	136	SER	-	expression tag	UNP A4VT01

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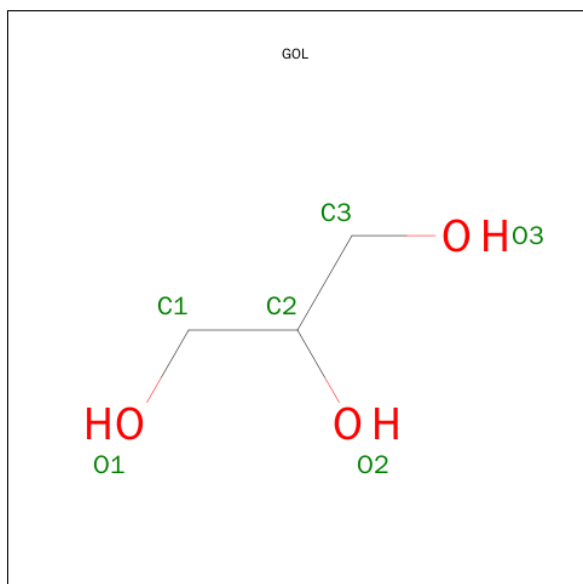
Chain	Residue	Modelled	Actual	Comment	Reference
C	137	HIS	-	expression tag	UNP A4VT01
C	138	MSE	-	expression tag	UNP A4VT01
D	118	MSE	-	expression tag	UNP A4VT01
D	119	GLY	-	expression tag	UNP A4VT01
D	120	SER	-	expression tag	UNP A4VT01
D	121	SER	-	expression tag	UNP A4VT01
D	122	HIS	-	expression tag	UNP A4VT01
D	123	HIS	-	expression tag	UNP A4VT01
D	124	HIS	-	expression tag	UNP A4VT01
D	125	HIS	-	expression tag	UNP A4VT01
D	126	HIS	-	expression tag	UNP A4VT01
D	127	HIS	-	expression tag	UNP A4VT01
D	128	SER	-	expression tag	UNP A4VT01
D	129	SER	-	expression tag	UNP A4VT01
D	130	GLY	-	expression tag	UNP A4VT01
D	131	LEU	-	expression tag	UNP A4VT01
D	132	VAL	-	expression tag	UNP A4VT01
D	133	PRO	-	expression tag	UNP A4VT01
D	134	ARG	-	expression tag	UNP A4VT01
D	135	GLY	-	expression tag	UNP A4VT01
D	136	SER	-	expression tag	UNP A4VT01
D	137	HIS	-	expression tag	UNP A4VT01
D	138	MSE	-	expression tag	UNP A4VT01
E	118	MSE	-	expression tag	UNP A4VT01
E	119	GLY	-	expression tag	UNP A4VT01
E	120	SER	-	expression tag	UNP A4VT01
E	121	SER	-	expression tag	UNP A4VT01
E	122	HIS	-	expression tag	UNP A4VT01
E	123	HIS	-	expression tag	UNP A4VT01
E	124	HIS	-	expression tag	UNP A4VT01
E	125	HIS	-	expression tag	UNP A4VT01
E	126	HIS	-	expression tag	UNP A4VT01
E	127	HIS	-	expression tag	UNP A4VT01
E	128	SER	-	expression tag	UNP A4VT01
E	129	SER	-	expression tag	UNP A4VT01
E	130	GLY	-	expression tag	UNP A4VT01
E	131	LEU	-	expression tag	UNP A4VT01
E	132	VAL	-	expression tag	UNP A4VT01
E	133	PRO	-	expression tag	UNP A4VT01
E	134	ARG	-	expression tag	UNP A4VT01
E	135	GLY	-	expression tag	UNP A4VT01
E	136	SER	-	expression tag	UNP A4VT01

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Chain	Residue	Modelled	Actual	Comment	Reference
E	137	HIS	-	expression tag	UNP A4VT01
E	138	MSE	-	expression tag	UNP A4VT01

- Molecule 2 is GLYCEROL (three-letter code: GOL) (formula:  $C_3H_8O_3$ ).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			6	3	3		
2	A	1	Total	C	O	0	0
			6	3	3		
2	B	1	Total	C	O	0	0
			6	3	3		
2	B	1	Total	C	O	0	0
			6	3	3		
2	B	1	Total	C	O	0	0
			6	3	3		
2	C	1	Total	C	O	0	0
			6	3	3		
2	D	1	Total	C	O	0	0
			6	3	3		
2	D	1	Total	C	O	0	0
			6	3	3		
2	E	1	Total	C	O	0	0
			6	3	3		
2	E	1	Total	C	O	0	0
			6	3	3		

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

- Molecule 3 is water.

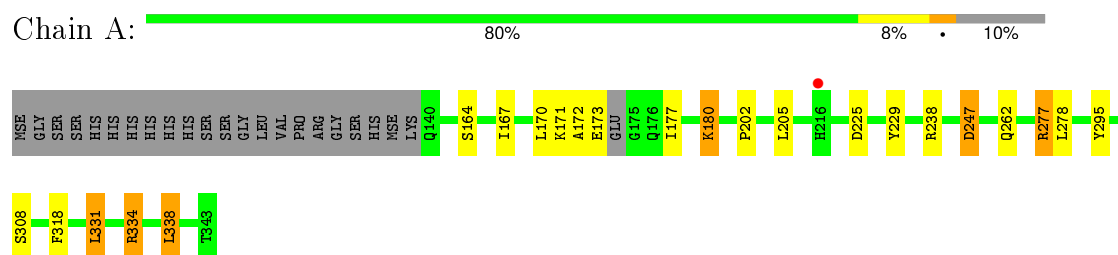
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	209	Total	O	0	0
			209	209		
3	B	183	Total	O	0	0
			183	183		
3	C	219	Total	O	0	0
			219	219		
3	D	227	Total	O	0	0
			227	227		
3	E	240	Total	O	0	0
			240	240		



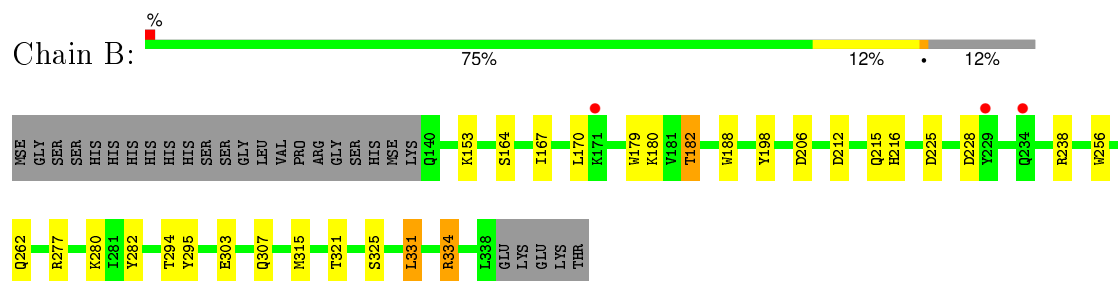
### 3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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 ( $\text{RSRZ} > 2$ ). 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.

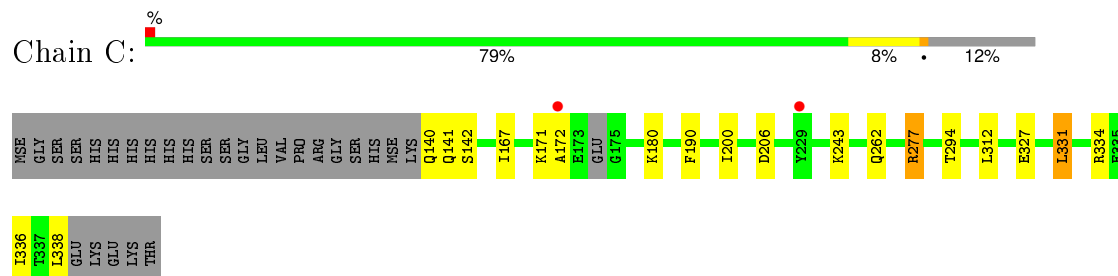
- Molecule 1: Translation initiation factor 2 (IF-2 GTPase)



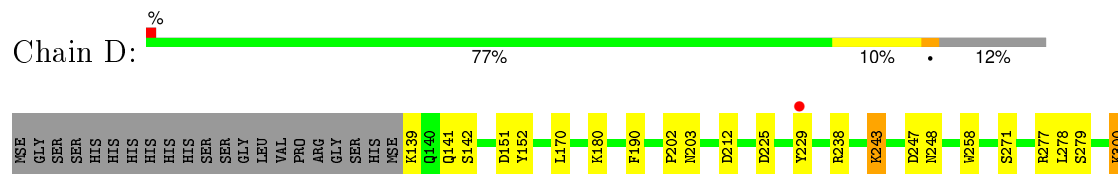
- Molecule 1: Translation initiation factor 2 (IF-2 GTPase)

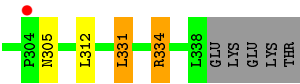


- Molecule 1: Translation initiation factor 2 (IF-2 GTPase)

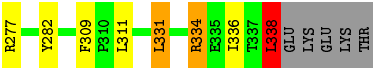
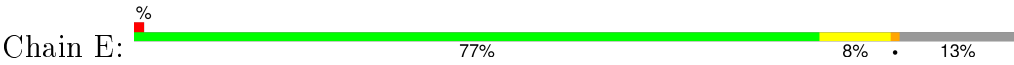


- Molecule 1: Translation initiation factor 2 (IF-2 GTPase)





● Molecule 1: Translation initiation factor 2 (IF-2 GTPase)



## 4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	127.08 Å   77.34 Å   131.64 Å 90.00°   115.93°   90.00°	Depositor
Resolution (Å)	39.46 – 1.50 39.46 – 1.50	Depositor EDS
% Data completeness (in resolution range)	97.0 (39.46-1.50) 97.0 (39.46-1.50)	Depositor EDS
$R_{merge}$	0.06	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.81 (at 1.50 Å)	Xtriage
Refinement program	REFMAC	Depositor
R, $R_{free}$	0.161 , 0.189 0.161 , 0.189	Depositor DCC
$R_{free}$ test set	8926 reflections (5.28%)	DCC
Wilson B-factor (Å <sup>2</sup> )	17.8	Xtriage
Anisotropy	0.296	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.38 , 44.6	EDS
Estimated twinning fraction	0.009 for h,-k,-h-l	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 178754 reflections	Xtriage
$F_o, F_c$ correlation	0.97	EDS
Total number of atoms	9071	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	20.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.97% 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, 0.375 respectively for untwinned datasets, and 0.333, 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	1.42	6/1645 (0.4%)	1.45	19/2231 (0.9%)
1	B	1.40	7/1612 (0.4%)	1.27	12/2187 (0.5%)
1	C	1.38	5/1614 (0.3%)	1.34	10/2188 (0.5%)
1	D	1.51	10/1629 (0.6%)	1.47	20/2210 (0.9%)
1	E	1.39	3/1605 (0.2%)	1.26	12/2176 (0.6%)
All	All	1.42	31/8105 (0.4%)	1.36	73/10992 (0.7%)

The worst 5 of 31 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	E	262	GLN	CB-CG	-11.85	1.20	1.52
1	A	262	GLN	CB-CG	-10.45	1.24	1.52
1	D	334	ARG	CB-CG	-8.99	1.28	1.52
1	C	262	GLN	CB-CG	-7.87	1.31	1.52
1	D	334	ARG	CD-NE	-7.66	1.33	1.46

The worst 5 of 73 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	277	ARG	NE-CZ-NH2	-27.96	106.32	120.30
1	D	334	ARG	NE-CZ-NH2	-18.30	111.15	120.30
1	C	277	ARG	NE-CZ-NH2	-17.26	111.67	120.30
1	A	277	ARG	NE-CZ-NH1	16.40	128.50	120.30
1	C	334	ARG	NE-CZ-NH2	-15.74	112.43	120.30

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	1606	0	1521	10	0
1	B	1579	0	1516	18	0
1	C	1579	0	1524	8	0
1	D	1590	0	1514	10	0
1	E	1573	0	1513	15	0
2	A	12	0	16	0	0
2	B	18	0	24	0	0
2	C	6	0	8	0	0
2	D	12	0	16	0	0
2	E	18	0	24	3	0
3	A	209	0	0	3	0
3	B	183	0	0	5	0
3	C	219	0	0	2	0
3	D	227	0	0	3	0
3	E	240	0	0	5	0
All	All	9071	0	7676	62	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.

The worst 5 of 62 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:C:243:LYS:HG2	3:C:513:HOH:O	1.57	1.05
1:D:142:SER:O	1:D:334:ARG:HD3	1.57	1.04
1:B:170:LEU:HD13	1:B:179:TRP:HA	1.57	0.84
1:D:142:SER:O	1:D:334:ARG:CD	2.27	0.82
1:A:247:ASP:O	3:A:502:HOH:O	1.96	0.81

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	A	200/226 (88%)	194 (97%)	6 (3%)	0	100	100
1	B	197/226 (87%)	191 (97%)	6 (3%)	0	100	100
1	C	195/226 (86%)	187 (96%)	8 (4%)	0	100	100
1	D	199/226 (88%)	193 (97%)	6 (3%)	0	100	100
1	E	193/226 (85%)	187 (97%)	6 (3%)	0	100	100
All	All	984/1130 (87%)	952 (97%)	32 (3%)	0	100	100

There are no Ramachandran outliers to report.

### 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	180/201 (90%)	177 (98%)	3 (2%)	68	37
1	B	177/201 (88%)	175 (99%)	2 (1%)	80	58
1	C	178/201 (89%)	174 (98%)	4 (2%)	60	25
1	D	179/201 (89%)	176 (98%)	3 (2%)	68	37
1	E	178/201 (89%)	175 (98%)	3 (2%)	68	37
All	All	892/1005 (89%)	877 (98%)	15 (2%)	68	37

5 of 15 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	C	171	LYS
1	C	277	ARG
1	E	171	LYS
1	C	167	ILE
1	D	331	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 7 such sidechains are listed below:

Mol	Chain	Res	Type
1	C	262	GLN
1	E	262	GLN
1	D	203	ASN
1	A	262	GLN
1	E	234	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](#)

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](#)

11 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
2	GOL	A	401	-	5,5,5	0.75	0	5,5,5	0.45	0
2	GOL	A	402	-	5,5,5	0.67	0	5,5,5	1.13	0
2	GOL	B	401	-	5,5,5	0.41	0	5,5,5	1.29	1 (20%)
2	GOL	B	402	-	5,5,5	0.52	0	5,5,5	0.73	0
2	GOL	B	403	-	5,5,5	0.33	0	5,5,5	0.85	0
2	GOL	C	401	-	5,5,5	0.41	0	5,5,5	1.27	1 (20%)
2	GOL	D	401	-	5,5,5	0.51	0	5,5,5	0.59	0
2	GOL	D	402	-	5,5,5	0.63	0	5,5,5	1.41	1 (20%)
2	GOL	E	401	-	5,5,5	0.78	0	5,5,5	0.84	0
2	GOL	E	402	-	5,5,5	0.54	0	5,5,5	1.23	1 (20%)
2	GOL	E	403	-	5,5,5	0.53	0	5,5,5	1.42	1 (20%)

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	GOL	A	401	-	-	0/4/4/4	0/0/0/0
2	GOL	A	402	-	-	0/4/4/4	0/0/0/0
2	GOL	B	401	-	-	0/4/4/4	0/0/0/0
2	GOL	B	402	-	-	0/4/4/4	0/0/0/0
2	GOL	B	403	-	-	0/4/4/4	0/0/0/0
2	GOL	C	401	-	-	0/4/4/4	0/0/0/0
2	GOL	D	401	-	-	0/4/4/4	0/0/0/0
2	GOL	D	402	-	-	0/4/4/4	0/0/0/0
2	GOL	E	401	-	-	0/4/4/4	0/0/0/0
2	GOL	E	402	-	-	0/4/4/4	0/0/0/0
2	GOL	E	403	-	-	0/4/4/4	0/0/0/0

There are no bond length outliers.

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	401	GOL	C3-C2-C1	-2.33	101.24	111.06
2	B	401	GOL	C3-C2-C1	-2.29	101.41	111.06
2	E	403	GOL	C3-C2-C1	-2.17	101.91	111.06
2	E	402	GOL	O3-C3-C2	-2.16	99.03	109.97
2	D	402	GOL	O2-C2-C1	2.64	121.17	108.47

There are no chirality outliers.



There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	E	403	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	202/226 (89%)	-0.28	1 (0%) 91 93	13, 18, 32, 46	0
1	B	198/226 (87%)	-0.20	3 (1%) 76 79	13, 19, 31, 37	0
1	C	197/226 (87%)	-0.18	2 (1%) 84 86	12, 17, 30, 41	0
1	D	199/226 (88%)	-0.12	2 (1%) 84 86	11, 16, 30, 38	0
1	E	195/226 (86%)	-0.31	3 (1%) 76 79	11, 16, 27, 36	0
All	All	991/1130 (87%)	-0.22	11 (1%) 82 85	11, 17, 30, 46	0

The worst 5 of 11 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	229	TYR	3.6
1	C	172	ALA	3.4
1	D	304	PRO	3.4
1	C	229	TYR	3.3
1	B	234	GLN	2.7

### 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. LLDF column lists the quality of electron

density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors(Å <sup>2</sup> )	Q<0.9
2	GOL	B	401	6/6	0.91	0.17	7.19	24,30,35,40	0
2	GOL	E	403	6/6	0.95	0.11	4.96	20,26,33,41	0
2	GOL	D	402	6/6	0.92	0.13	3.08	18,23,32,33	0
2	GOL	D	401	6/6	0.98	0.10	2.94	13,14,16,16	0
2	GOL	A	402	6/6	0.96	0.08	2.84	18,22,26,34	0
2	GOL	C	401	6/6	0.90	0.11	2.77	21,25,30,38	0
2	GOL	E	402	6/6	0.95	0.08	2.27	16,19,25,34	0
2	GOL	B	403	6/6	0.98	0.13	1.99	18,24,26,33	0
2	GOL	E	401	6/6	0.98	0.07	1.15	12,14,15,16	0
2	GOL	A	401	6/6	0.98	0.07	0.46	14,15,15,16	0
2	GOL	B	402	6/6	0.98	0.07	-0.56	15,17,18,20	0

## 6.5 Other polymers

There are no such residues in this entry.