



Full wwPDB X-ray Structure Validation Report ⓘ

Nov 14, 2023 – 02:24 PM JST

PDB ID : 5ZH8
Title : Crystal Structure of FmtA from Staphylococcus aureus at 2.58 Å
Authors : Dalal, V.; Kumar, P.; Golemi-Kotra, D.; Kumar, P.
Deposited on : 2018-03-12
Resolution : 2.60 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

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>.

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 : 2.36
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) : 2.36

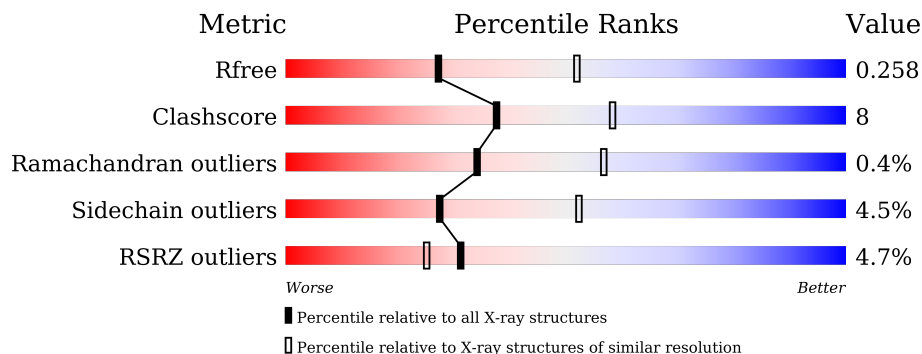
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.60 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	3163 (2.60-2.60)
Clashscore	141614	3518 (2.60-2.60)
Ramachandran outliers	138981	3455 (2.60-2.60)
Sidechain outliers	138945	3455 (2.60-2.60)
RSRZ outliers	127900	3104 (2.60-2.60)

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 ≥ 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	397	
1	B	397	

2 Entry composition [i](#)

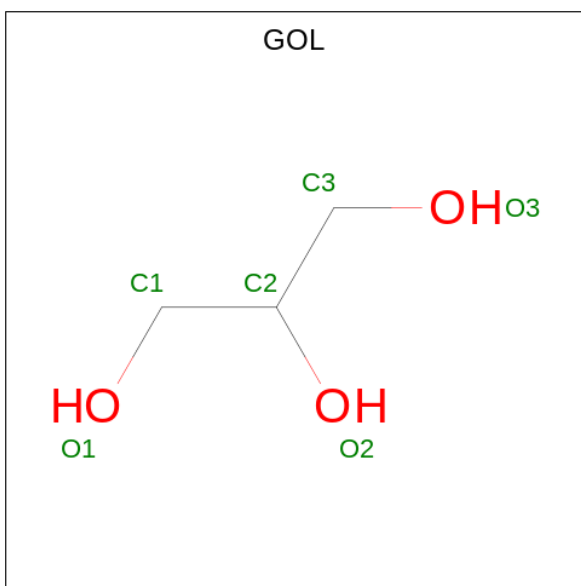
There are 3 unique types of molecules in this entry. The entry contains 5927 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Protein FmtA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	358	Total	C	N	O	S	0	0	0
			2928	1902	483	537	6			
1	B	349	Total	C	N	O	S	0	0	0
			2851	1851	470	524	6			

- Molecule 2 is GLYCEROL (three-letter code: GOL) (formula: C₃H₈O₃).



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

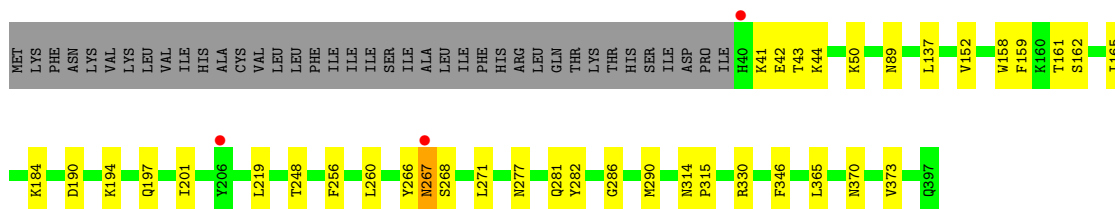
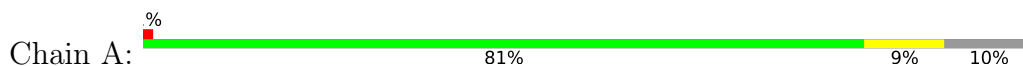
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	72	Total 72	O 72	0	0
3	B	64	Total 64	O 64	0	0

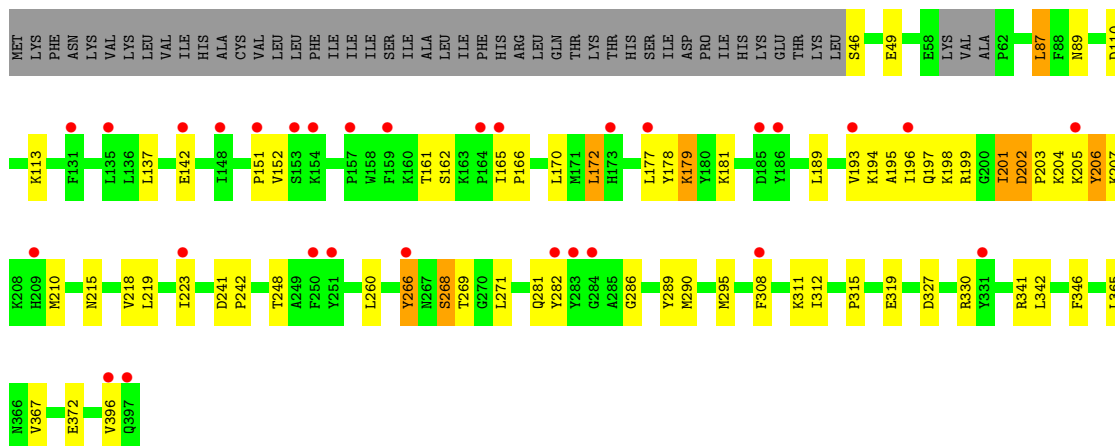
3 Residue-property plots [i](#)

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: Protein FmtA



- Molecule 1: Protein FmtA



4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	105.39Å 128.23Å 107.60Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	50.00 – 2.60 47.32 – 2.60	Depositor EDS
% Data completeness (in resolution range)	97.6 (50.00-2.60) 97.6 (47.32-2.60)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.60 (at 2.61Å)	Xtrriage
Refinement program	REFMAC 5.8.0216	Depositor
R, R_{free}	0.208 , 0.254 0.215 , 0.258	Depositor DCC
R_{free} test set	1118 reflections (5.03%)	wwPDB-VP
Wilson B-factor (Å ²)	53.7	Xtrriage
Anisotropy	0.354	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 55.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.34$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	5927	wwPDB-VP
Average B, all atoms (Å ²)	72.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.79% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical 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: 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	0.34	0/3000	0.52	0/4040
1	B	0.36	0/2921	0.55	0/3934
All	All	0.35	0/5921	0.53	0/7974

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	2928	0	2952	21	0
1	B	2851	0	2862	71	0
2	A	6	0	8	0	0
2	B	6	0	8	0	0
3	A	72	0	0	0	0
3	B	64	0	0	0	0
All	All	5927	0	5830	89	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 8.

All (89) 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:B:178:TYR:CZ	1:B:202:ASP:HB2	1.59	1.38
1:B:266:TYR:CE2	1:B:271:LEU:HD13	1.63	1.32
1:B:178:TYR:CE1	1:B:202:ASP:HB2	1.82	1.13
1:B:142:GLU:OE2	1:B:312:ILE:HG21	1.47	1.13
1:B:266:TYR:CZ	1:B:271:LEU:HD13	1.96	0.99
1:B:202:ASP:OD2	1:B:205:LYS:HB2	1.67	0.95
1:B:266:TYR:CE2	1:B:271:LEU:CD1	2.54	0.91
1:B:178:TYR:CZ	1:B:202:ASP:CB	2.54	0.87
1:B:178:TYR:CE2	1:B:202:ASP:HB2	2.14	0.82
1:B:178:TYR:CE1	1:B:202:ASP:CB	2.64	0.79
1:B:202:ASP:OD2	1:B:205:LYS:CB	2.32	0.77
1:B:215:ASN:O	1:B:219:LEU:HD12	1.85	0.76
1:B:202:ASP:OD2	1:B:205:LYS:CG	2.34	0.75
1:B:152:VAL:HG21	1:B:170:LEU:HD11	1.69	0.74
1:B:193:VAL:HA	1:B:196:ILE:HG22	1.68	0.73
1:B:199:ARG:O	1:B:199:ARG:HG2	1.89	0.73
1:A:184:LYS:HE2	1:A:277:ASN:HD22	1.55	0.71
1:B:268:SER:O	1:B:269:THR:OG1	2.09	0.70
1:B:179:LYS:HB2	1:B:210:MET:HE1	1.73	0.70
1:A:370:ASN:HD21	1:B:266:TYR:HE1	1.41	0.68
1:B:266:TYR:OH	1:B:367:VAL:HG11	1.93	0.67
1:A:184:LYS:CE	1:A:277:ASN:HD22	2.08	0.66
1:B:110:ASP:OD2	1:B:113:LYS:HG3	1.98	0.64
1:B:202:ASP:OD2	1:B:205:LYS:HD2	1.98	0.63
1:B:199:ARG:HB3	1:B:199:ARG:HH11	1.64	0.62
1:B:142:GLU:OE1	1:B:308:PHE:CE1	2.52	0.62
1:B:266:TYR:CZ	1:B:271:LEU:CD1	2.80	0.62
1:A:165:ILE:HD11	1:A:201:ILE:HD11	1.81	0.61
1:A:267:ASN:OD1	1:A:268:SER:N	2.34	0.60
1:B:152:VAL:CG2	1:B:170:LEU:HD11	2.32	0.59
1:B:202:ASP:OD2	1:B:205:LYS:CD	2.51	0.59
1:B:194:LYS:O	1:B:197:GLN:HG3	2.01	0.59
1:B:202:ASP:OD1	1:B:203:PRO:O	2.23	0.57
1:A:267:ASN:OD1	1:A:267:ASN:C	2.44	0.56
1:B:205:LYS:HE2	1:B:210:MET:SD	2.47	0.55
1:B:179:LYS:HB2	1:B:210:MET:CE	2.36	0.54
1:B:142:GLU:OE1	1:B:308:PHE:CD1	2.61	0.54
1:B:248:THR:HG23	1:B:290:MET:HG2	1.90	0.54
1:A:42:GLU:HG2	1:A:43:THR:N	2.23	0.53
1:A:248:THR:HG23	1:A:290:MET:HG2	1.90	0.52
1:B:195:ALA:O	1:B:198:LYS:HB2	2.10	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:170:LEU:HB3	1:B:219:LEU:HD11	1.91	0.52
1:B:179:LYS:CB	1:B:210:MET:HE1	2.38	0.52
1:A:43:THR:HG23	1:A:43:THR:O	2.09	0.52
1:B:201:ILE:HD11	1:B:206:TYR:HA	1.91	0.52
1:B:170:LEU:CB	1:B:219:LEU:HD11	2.41	0.51
1:B:266:TYR:CD1	1:B:266:TYR:N	2.73	0.51
1:B:199:ARG:NH1	1:B:199:ARG:CB	2.73	0.51
1:A:314:ASN:HB2	1:A:315:PRO:HD3	1.94	0.50
1:A:184:LYS:CE	1:A:277:ASN:ND2	2.75	0.50
1:B:165:ILE:HD12	1:B:165:ILE:O	2.10	0.50
1:B:295:MET:CE	1:B:342:LEU:HD22	2.42	0.50
1:A:184:LYS:HE2	1:A:277:ASN:ND2	2.24	0.50
1:B:189:LEU:O	1:B:193:VAL:HG22	2.12	0.49
1:B:295:MET:HE1	1:B:342:LEU:HD22	1.95	0.49
1:B:319:GLU:HB2	1:B:396:VAL:HG22	1.95	0.49
1:A:370:ASN:ND2	1:B:266:TYR:HE1	2.11	0.48
1:A:42:GLU:HG3	1:A:44:LYS:HG2	1.96	0.47
1:A:256:PHE:O	1:A:260:LEU:HD13	2.14	0.47
1:B:199:ARG:HB3	1:B:199:ARG:NH1	2.26	0.47
1:A:137:LEU:HD11	1:A:219:LEU:HD13	1.97	0.47
1:A:159:PHE:HA	1:A:197:GLN:HG3	1.96	0.46
1:B:137:LEU:CD2	1:B:223:ILE:HD11	2.45	0.46
1:A:158:TRP:CZ3	1:A:194:LYS:HG3	2.50	0.46
1:B:202:ASP:OD2	1:B:205:LYS:HG3	2.15	0.45
1:B:266:TYR:CD2	1:B:271:LEU:HB2	2.51	0.45
1:B:266:TYR:CD2	1:B:271:LEU:HD13	2.35	0.45
1:B:266:TYR:CD2	1:B:271:LEU:CD1	2.96	0.45
1:B:315:PRO:HA	1:B:396:VAL:CG2	2.46	0.45
1:B:201:ILE:HD13	1:B:202:ASP:N	2.32	0.45
1:B:137:LEU:HD23	1:B:223:ILE:HD11	1.99	0.44
1:B:46:SER:HB2	1:B:49:GLU:HG2	1.98	0.44
1:B:179:LYS:HB2	1:B:179:LYS:HE2	1.42	0.44
1:B:172:LEU:HD11	1:B:327:ASP:HB2	2.00	0.43
1:B:89:ASN:O	1:B:365:LEU:HD12	2.18	0.43
1:B:202:ASP:OD1	1:B:202:ASP:C	2.56	0.43
1:B:170:LEU:HD13	1:B:218:VAL:HG13	2.00	0.43
1:A:89:ASN:O	1:A:365:LEU:HD12	2.19	0.43
1:A:266:TYR:CZ	1:A:271:LEU:HD13	2.54	0.42
1:B:199:ARG:HH11	1:B:199:ARG:CB	2.28	0.42
1:A:373:VAL:HG12	1:B:87:LEU:HD23	2.02	0.42
1:B:203:PRO:O	1:B:204:LYS:CB	2.67	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:260:LEU:HD22	1:B:289:TYR:CG	2.55	0.41
1:B:142:GLU:OE2	1:B:312:ILE:CG2	2.40	0.41
1:B:170:LEU:HB3	1:B:219:LEU:CD1	2.51	0.41
1:B:241:ASP:HB2	1:B:242:PRO:HD3	2.03	0.41
1:B:341:ARG:NH1	1:B:372:GLU:OE2	2.54	0.40
1:B:151:PRO:HA	1:B:166:PRO:HA	2.03	0.40
1:B:193:VAL:HA	1:B:196:ILE:CG2	2.45	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	356/397 (90%)	340 (96%)	15 (4%)	1 (0%)	41	64
1	B	345/397 (87%)	327 (95%)	16 (5%)	2 (1%)	25	47
All	All	701/794 (88%)	667 (95%)	31 (4%)	3 (0%)	34	57

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	206	TYR
1	A	286	GLY
1	B	286	GLY

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	317/358 (88%)	306 (96%)	11 (4%)	36	62
1	B	309/358 (86%)	292 (94%)	17 (6%)	21	43
All	All	626/716 (87%)	598 (96%)	28 (4%)	27	52

All (28) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	41	LYS
1	A	50	LYS
1	A	152	VAL
1	A	161	THR
1	A	162	SER
1	A	190	ASP
1	A	267	ASN
1	A	281	GLN
1	A	282	TYR
1	A	330	ARG
1	A	346	PHE
1	B	87	LEU
1	B	161	THR
1	B	162	SER
1	B	172	LEU
1	B	177	LEU
1	B	179	LYS
1	B	181	LYS
1	B	201	ILE
1	B	202	ASP
1	B	207	LYS
1	B	266	TYR
1	B	268	SER
1	B	281	GLN
1	B	282	TYR
1	B	311	LYS
1	B	330	ARG
1	B	346	PHE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	144	HIS
1	A	234	ASN
1	A	254	GLN
1	A	318	HIS
1	B	314	ASN
1	B	343	ASN

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 monosaccharides in this entry.

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	$\# Z > 2$
2	GOL	A	401	-	5,5,5	0.26	0	5,5,5	0.22	0
2	GOL	B	401	-	5,5,5	0.31	0	5,5,5	0.17	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. '2' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	GOL	A	401	-	-	2/4/4/4	-
2	GOL	B	401	-	-	3/4/4/4	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (5) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	401	GOL	O1-C1-C2-C3
2	B	401	GOL	O2-C2-C3-O3
2	B	401	GOL	C1-C2-C3-O3
2	A	401	GOL	O1-C1-C2-O2
2	B	401	GOL	O1-C1-C2-C3

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

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, 95th 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 ²)	Q<0.9
1	A	358/397 (90%)	-0.21	3 (0%) 86 84	37, 55, 93, 139	0
1	B	349/397 (87%)	0.45	30 (8%) 10 7	45, 84, 142, 161	0
All	All	707/794 (89%)	0.12	33 (4%) 31 25	37, 65, 128, 161	0

All (33) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	266	TYR	5.1
1	B	142	GLU	4.7
1	B	148	ILE	4.3
1	B	165	ILE	4.2
1	B	151	PRO	4.2
1	B	159	PHE	4.1
1	B	223	ILE	3.6
1	B	209	HIS	3.5
1	A	40	HIS	3.4
1	B	205	LYS	3.4
1	B	153	SER	3.3
1	B	308	PHE	3.2
1	B	193	VAL	3.2
1	A	267	ASN	3.2
1	B	157	PRO	3.0
1	B	282	TYR	2.9
1	A	206	TYR	2.9
1	B	185	ASP	2.8
1	B	177	LEU	2.8
1	B	196	ILE	2.6
1	B	250	PHE	2.6
1	B	164	PRO	2.5
1	B	173	HIS	2.5
1	B	397	GLN	2.4

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Mol	Chain	Res	Type	RSRZ
1	B	251	TYR	2.3
1	B	283	TYR	2.3
1	B	154	LYS	2.3
1	B	284	GLY	2.2
1	B	186	TYR	2.2
1	B	131	PHE	2.1
1	B	135	LEU	2.0
1	B	396	VAL	2.0
1	B	331	TYR	2.0

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 monosaccharides 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, 95th 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 ²)	Q<0.9
2	GOL	B	401	6/6	0.68	0.29	89,91,91,91	0
2	GOL	A	401	6/6	0.73	0.30	85,86,87,88	0

6.5 Other polymers [i](#)

There are no such residues in this entry.