



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 13, 2017 – 11:58 pm GMT

PDB ID : 3TFH  
Title : DMSP-dependent demethylase from *P. ubiquus* - apo  
Authors : Schuller, D.J.; Reisch, C.R.; Moran, M.A.; Whitman, W.B.; Lanzilotta, W.N.  
Deposited on : 2011-08-15  
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

<http://wwpdb.org/validation/2016/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

---

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467  
Mogul : 1.7.2 (RC1), CSD as538be (2017)  
Xtriage (Phenix) : 1.9-1692  
EDS : trunk28620  
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)  
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) : recalc28949

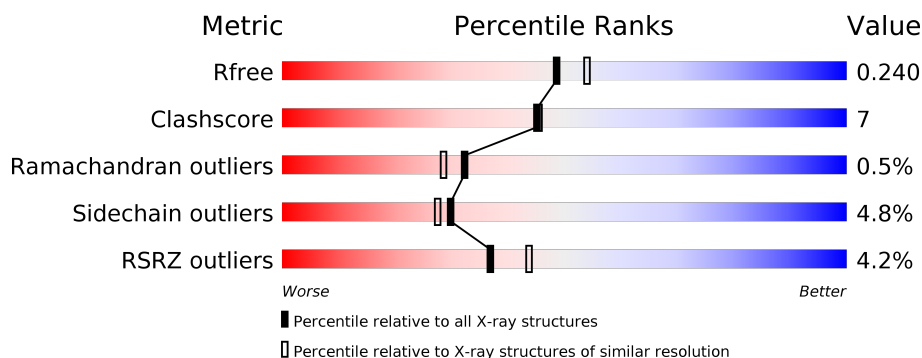
# 1 Overall quality at a glance

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))
$R_{free}$	100719	4243 (2.10-2.10)
Clashscore	112137	4788 (2.10-2.10)
Ramachandran outliers	110173	4740 (2.10-2.10)
Sidechain outliers	110143	4741 (2.10-2.10)
RSRZ outliers	101464	4275 (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.

Mol	Chain	Length	Quality of chain
1	A	369	<div> <div>4%</div> <div> <div></div> <div>83%</div> <div>15%</div> <div></div> </div> </div>
1	B	369	<div> <div>5%</div> <div> <div></div> <div>85%</div> <div>13%</div> <div></div> </div> </div>

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	370	-	-	-	X
2	GOL	A	371	-	-	-	X

## 2 Entry composition [i](#)

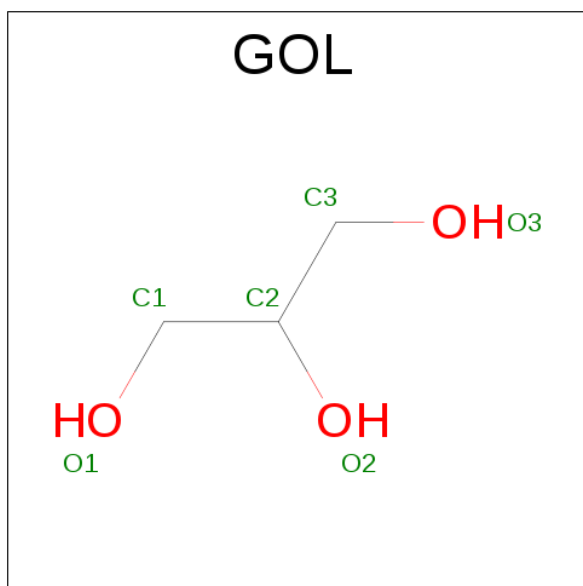
There are 4 unique types of molecules in this entry. The entry contains 6668 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 GcvT-like Aminomethyltransferase protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	369	Total	C	N	O	S	0	25	0
			3091	1996	509	573	13			
1	B	369	Total	C	N	O	S	0	10	0
			3004	1926	499	565	14			

- Molecule 2 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
2	A	1	Total	C	O	0	0
			6	3	3		
2	A	1	Total	C	O	0	0
			6	3	3		

- Molecule 3 is SODIUM ION (three-letter code: NA) (formula: Na).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	1	Total 1	Na 1	0	0
3	A	1	Total 1	Na 1	0	0

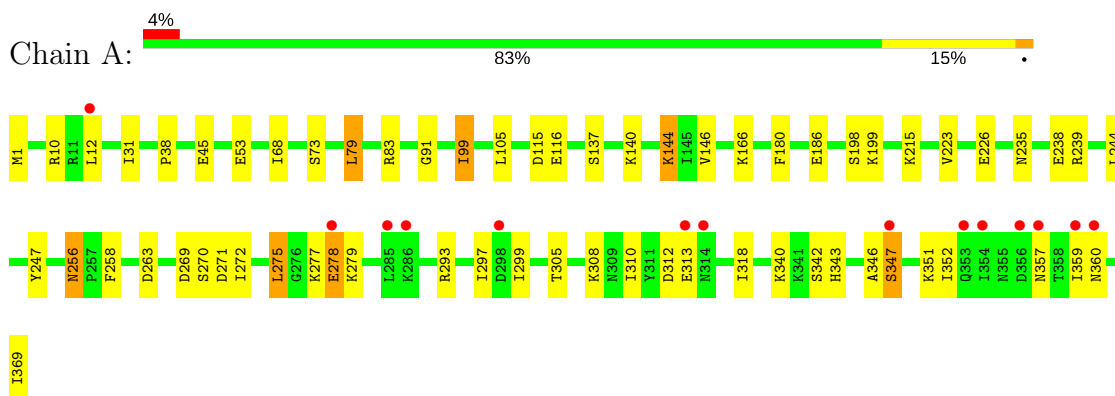
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	256	Total 256	O 256	0	0
4	B	303	Total 303	O 303	0	0

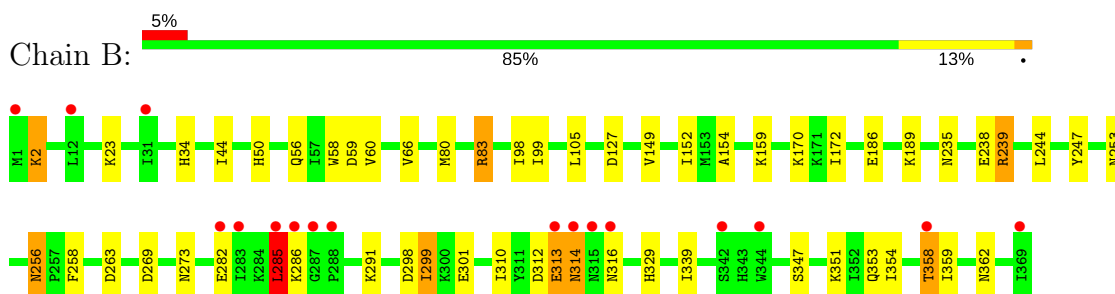
### 3 Residue-property plots [i](#)

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: GcvT-like Aminomethyltransferase protein



#### • Molecule 1: GcvT-like Aminomethyltransferase protein



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	53.66Å 121.73Å 59.65Å 90.00° 101.01° 90.00°	Depositor
Resolution (Å)	43.51 – 2.10 43.51 – 2.10	Depositor EDS
% Data completeness (in resolution range)	100.0 (43.51-2.10) 100.0 (43.51-2.10)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	0.07	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	15.04 (at 2.10Å)	Xtriage
Refinement program	REFMAC 5.5.0109	Depositor
R, $R_{free}$	0.166 , 0.231 0.175 , 0.240	Depositor DCC
$R_{free}$ test set	1066 reflections (2.49%)	DCC
Wilson B-factor (Å <sup>2</sup> )	26.1	Xtriage
Anisotropy	0.298	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.34 , 55.6	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.96	EDS
Total number of atoms	6668	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	43.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.39% 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, NA

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	0.95	0/3228	0.87	4/4334 (0.1%)
1	B	0.99	2/3092 (0.1%)	0.85	3/4159 (0.1%)
All	All	0.97	2/6320 (0.0%)	0.86	7/8493 (0.1%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	66	VAL	CB-CG1	6.02	1.65	1.52
1	B	149	VAL	CB-CG1	6.00	1.65	1.52

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	293	ARG	NE-CZ-NH2	-6.73	116.94	120.30
1	B	285	LEU	CA-CB-CG	6.22	129.62	115.30
1	A	83	ARG	NE-CZ-NH1	5.98	123.29	120.30
1	A	293	ARG	NE-CZ-NH1	5.82	123.21	120.30
1	A	10	ARG	NE-CZ-NH1	5.35	122.97	120.30
1	B	59	ASP	CB-CG-OD2	5.16	122.94	118.30
1	B	83	ARG	NE-CZ-NH2	-5.03	117.79	120.30

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	A	3091	0	3164	61	1
1	B	3004	0	3009	30	1
2	A	12	0	16	2	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
4	A	256	0	0	9	0
4	B	303	0	0	9	0
All	All	6668	0	6189	91	1

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.

All (91) 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:A:305:THR:HG21	4:B:380:HOH:O	1.56	1.04
1:A:299:ILE:HG23	1:A:359:ILE:CD1	1.90	1.01
1:A:305:THR:CG2	4:B:380:HOH:O	2.12	0.94
1:A:99:ILE:C	1:A:99:ILE:HD12	1.90	0.92
1:A:312:ASP:O	4:A:460:HOH:O	1.93	0.87
1:A:79:LEU:HD22	1:A:275:LEU:HD13	1.57	0.85
1:A:186[B]:GLU:OE1	1:A:215[B]:LYS:HE2	1.81	0.81
1:A:278[A]:GLU:OE1	1:A:279:LYS:N	2.14	0.79
1:A:140[B]:LYS:HE3	1:A:140[B]:LYS:HA	1.64	0.79
1:A:140[B]:LYS:HA	1:A:140[B]:LYS:CE	2.16	0.76
1:A:235:ASN:HD22	1:A:238:GLU:H	1.37	0.73
1:A:299:ILE:CG2	1:A:359:ILE:CD1	2.68	0.71
1:A:299:ILE:HG23	1:A:359:ILE:HD11	1.73	0.71
1:B:186:GLU:OE1	4:B:556:HOH:O	2.10	0.70
1:A:199:LYS:HE3	1:A:244:LEU:HD21	1.73	0.70
1:B:310:ILE:HD11	1:B:339:ILE:CD1	2.22	0.69
1:A:99:ILE:O	1:A:99:ILE:HD12	1.93	0.69
1:A:215[B]:LYS:CE	4:A:525:HOH:O	2.40	0.69
1:B:354:ILE:HD12	1:B:359:ILE:HD12	1.75	0.69
1:A:99:ILE:HD11	1:A:275:LEU:O	1.92	0.68
1:A:297:ILE:HG22	1:A:299:ILE:HG12	1.75	0.67
1:A:79:LEU:HD22	1:A:275:LEU:CD1	2.24	0.67
1:B:235:ASN:HD22	1:B:238:GLU:H	1.44	0.66
1:A:99:ILE:C	1:A:99:ILE:CD1	2.64	0.66
1:B:282:GLU:O	1:B:285:LEU:HD12	1.95	0.66

*Continued on next page...*

*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:189:LYS:HE3	4:B:538:HOH:O	1.96	0.65
1:A:299:ILE:HG23	1:A:359:ILE:HD13	1.77	0.64
1:A:45:GLU:HG3	4:A:387:HOH:O	1.97	0.64
1:A:256:ASN:HD22	1:A:258:PHE:H	1.44	0.64
1:A:215[B]:LYS:HE3	4:A:525:HOH:O	1.98	0.63
1:A:272:ILE:O	1:A:277:LYS:NZ	2.23	0.62
1:A:369:ILE:HD12	1:A:369:ILE:C	2.20	0.62
1:B:310:ILE:HD11	1:B:339:ILE:HD13	1.83	0.61
1:A:12[B]:LEU:HD11	1:A:31:ILE:HG13	1.84	0.59
1:B:239:ARG:NE	4:B:605:HOH:O	2.04	0.59
1:A:312:ASP:HB3	1:A:318:ILE:HG21	1.85	0.58
1:A:166:LYS:HE2	1:A:223:VAL:HB	1.86	0.57
1:A:1:MET:HE2	4:A:419:HOH:O	2.04	0.56
1:B:273:ASN:ND2	4:B:514:HOH:O	2.40	0.54
1:A:299:ILE:CG2	1:A:359:ILE:HD13	2.37	0.54
1:B:50:HIS:HD2	1:B:56:GLN:HE21	1.55	0.54
1:A:342:SER:OG	1:A:343:HIS:HD2	1.91	0.53
1:B:312:ASP:OD1	1:B:314:ASN:ND2	2.42	0.52
1:A:247:TYR:OH	2:A:370:GOL:H11	2.10	0.52
1:A:12[B]:LEU:HG	1:A:31:ILE:CD1	2.40	0.51
1:A:38:PRO:HG2	4:A:446:HOH:O	2.10	0.51
1:A:115[B]:ASP:OD1	1:A:116:GLU:N	2.44	0.51
1:B:285:LEU:HD12	1:B:286:LYS:H	1.75	0.50
1:A:198:SER:HB2	1:A:238:GLU:OE1	2.12	0.49
1:A:369:ILE:HD12	1:A:369:ILE:O	2.13	0.49
1:B:50:HIS:CD2	1:B:56:GLN:HE21	2.30	0.49
1:A:215[B]:LYS:HE2	4:A:525:HOH:O	2.08	0.48
1:A:239:ARG:HB3	4:A:541:HOH:O	2.13	0.48
1:B:354:ILE:CD1	1:B:359:ILE:HD12	2.43	0.47
1:A:144[A]:LYS:CE	1:A:146:VAL:HG22	2.45	0.47
1:A:256:ASN:ND2	1:A:258:PHE:H	2.13	0.46
2:A:371:GOL:H31	1:B:247:TYR:OH	2.15	0.45
1:B:239:ARG:HD2	1:B:244:LEU:O	2.16	0.45
1:B:58:TRP:HB2	1:B:154:ALA:HB3	1.98	0.45
1:B:105:LEU:O	1:B:253:ASN:HA	2.17	0.45
1:B:256:ASN:HD22	1:B:258:PHE:H	1.64	0.45
1:A:312:ASP:HB3	1:A:318:ILE:CG2	2.46	0.45
1:A:369:ILE:CD1	1:A:369:ILE:C	2.85	0.44
1:A:308[B]:LYS:HE2	1:A:308[B]:LYS:HB2	1.67	0.44
1:A:271:ASP:O	4:A:563:HOH:O	2.21	0.43
1:A:318:ILE:HG21	1:A:343:HIS:CD2	2.53	0.43

*Continued on next page...*

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:256:ASN:C	1:B:256:ASN:HD22	2.21	0.43
1:B:159:LYS:NZ	4:B:659:HOH:O	2.51	0.43
1:A:144[A]:LYS:HE2	1:A:146:VAL:HG22	1.99	0.43
1:B:314:ASN:C	1:B:314:ASN:HD22	2.21	0.43
1:B:34:HIS:HE1	1:B:127:ASP:OD2	2.00	0.43
1:A:340[B]:LYS:HB3	1:A:340[B]:LYS:HE2	1.91	0.43
1:B:299:ILE:HD13	1:B:354:ILE:HD13	2.00	0.43
1:B:353:GLN:NE2	1:B:358:THR:OG1	2.52	0.42
1:A:91:GLY:HA2	1:A:180:PHE:CE2	2.55	0.42
1:B:80:MET:HB3	1:B:98:ILE:HD13	2.01	0.42
1:A:137:SER:O	1:A:140[B]:LYS:HE3	2.19	0.42
1:B:23[A]:LYS:CE	4:B:476:HOH:O	2.67	0.42
1:B:60:VAL:CG2	1:B:152:ILE:HG22	2.50	0.42
1:A:68:ILE:HG22	1:A:73:SER:HB2	2.02	0.42
1:A:12[B]:LEU:HG	1:A:31:ILE:HD12	2.02	0.41
1:A:351:LYS:NZ	1:A:360[A]:ASN:HD21	2.19	0.41
1:A:256:ASN:C	1:A:256:ASN:HD22	2.24	0.41
1:A:310:ILE:CG2	1:A:352:ILE:HD12	2.50	0.41
1:A:256:ASN:HD22	1:A:258:PHE:N	2.16	0.41
1:A:310:ILE:HG22	1:A:352:ILE:HD12	2.03	0.41
1:B:44:ILE:HD12	4:B:536:HOH:O	2.20	0.41
1:A:318:ILE:HG21	1:A:343:HIS:NE2	2.35	0.40
1:A:346:ALA:O	1:A:347:SER:HB2	2.22	0.40
1:B:99:ILE:HD12	1:B:99:ILE:HG23	1.90	0.40

All (1) 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:226:GLU:OE2	1:B:329:HIS:ND1[2_655]	2.18	0.02

## 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	392/369 (106%)	378 (96%)	13 (3%)	1 (0%)	44	44
1	B	376/369 (102%)	363 (96%)	10 (3%)	3 (1%)	22	17
All	All	768/738 (104%)	741 (96%)	23 (3%)	4 (0%)	32	28

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	347	SER
1	B	313	GLU
1	B	2	LYS
1	B	347	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	A	349/324 (108%)	335 (96%)	14 (4%)	36	36
1	B	334/324 (103%)	314 (94%)	20 (6%)	22	19
All	All	683/648 (105%)	649 (95%)	34 (5%)	30	26

All (34) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	79	LEU
1	A	99	ILE
1	A	105	LEU
1	A	144[A]	LYS
1	A	144[B]	LYS
1	A	256	ASN
1	A	263	ASP
1	A	269	ASP
1	A	270	SER
1	A	275	LEU

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Res	Type
1	A	278[A]	GLU
1	A	278[B]	GLU
1	A	313	GLU
1	A	357	ASN
1	B	2	LYS
1	B	83	ARG
1	B	170	LYS
1	B	172	ILE
1	B	239	ARG
1	B	256	ASN
1	B	263	ASP
1	B	269	ASP
1	B	285	LEU
1	B	291	LYS
1	B	298[A]	ASP
1	B	298[B]	ASP
1	B	299	ILE
1	B	301	GLU
1	B	313	GLU
1	B	314	ASN
1	B	316	ASN
1	B	351	LYS
1	B	358	THR
1	B	362	ASN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (20) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	65	GLN
1	A	228	ASN
1	A	235	ASN
1	A	253	ASN
1	A	256	ASN
1	A	314	ASN
1	A	343	HIS
1	B	24	GLN
1	B	34	HIS
1	B	50	HIS
1	B	54	HIS
1	B	156	GLN
1	B	228	ASN
1	B	235	ASN

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Res	Type
1	B	256	ASN
1	B	289	GLN
1	B	314	ASN
1	B	316	ASN
1	B	353	GLN
1	B	360	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 carbohydrates in this entry.

### 5.6 Ligand geometry [i](#)

Of 4 ligands modelled in this entry, 2 are monoatomic - leaving 2 for Mogul analysis.

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	370	-	5,5,5	0.41	0	5,5,5	0.60	0
2	GOL	A	371	-	5,5,5	0.47	0	5,5,5	1.26	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
2	GOL	A	370	-	-	0/4/4/4	0/0/0/0
2	GOL	A	371	-	-	0/4/4/4	0/0/0/0

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	370	GOL	1	0
2	A	371	GOL	1	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

### 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, 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	369/369 (100%)	0.05	14 (3%)	41	48	28, 40, 62, 78	0
1	B	369/369 (100%)	0.10	17 (4%)	33	39	25, 39, 66, 81	0
All	All	738/738 (100%)	0.08	31 (4%)	37	44	25, 39, 65, 81	0

All (31) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	285	LEU	6.5
1	B	1[A]	MET	5.9
1	A	356	ASP	4.5
1	A	313	GLU	4.1
1	B	286	LYS	3.6
1	A	353	GLN	3.1
1	A	359	ILE	3.1
1	B	315	ASN	3.0
1	A	298[A]	ASP	3.0
1	B	316	ASN	3.0
1	A	285	LEU	2.9
1	B	358	THR	2.8
1	B	313	GLU	2.8
1	A	360[A]	ASN	2.8
1	B	288	PRO	2.7
1	A	314	ASN	2.7
1	B	314	ASN	2.6
1	A	286[A]	LYS	2.6
1	B	342	SER	2.6
1	A	357	ASN	2.5
1	B	31	ILE	2.5
1	B	283	ILE	2.4
1	B	282	GLU	2.3
1	A	354	ILE	2.3

*Continued on next page...*



*Continued from previous page...*

Mol	Chain	Res	Type	RSRZ
1	B	344	TRP	2.3
1	A	12[A]	LEU	2.3
1	B	369	ILE	2.3
1	A	278[A]	GLU	2.2
1	A	347	SER	2.1
1	B	287	GLY	2.1
1	B	12	LEU	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 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	A	370	6/6	0.82	0.30	4.91	54,57,58,61	0
2	GOL	A	371	6/6	0.83	0.24	3.76	55,59,60,62	0
3	NA	B	370	1/1	0.99	0.19	0.95	34,34,34,34	0
3	NA	A	372	1/1	0.99	0.09	-2.03	33,33,33,33	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.