

# wwPDB X-ray Structure Validation Summary Report (i)

#### Aug 21, 2023 – 08:41 PM EDT

PDB ID	:	2OX4
Title	:	Crystal structure of putative dehydratase from Zymomonas mobilis ZM4
Authors	:	Patskovsky, Y.; Toro, R.; Sauder, J.M.; Freeman, J.C.; Bain, K.; Gheyi, T.;
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Deposited on	:	2007-02-19
Resolution	:	1.80  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 A user guide is available at https://www.wwpdb.org/validation/2017/XrayValidationReportHelp with specific help available everywhere you see the (i) 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 (1)) 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	:	2.35
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)	:	2.35

# 1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:  $X\text{-}RAY \, DIFFRACTION$ 

The reported resolution of this entry is 1.80 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.



Motria	Whole archive	Similar resolution
	$(\# { m Entries})$	$(\# { m Entries},  { m resolution}  { m range}({ m \AA}))$
R <sub>free</sub>	130704	5950 (1.80-1.80)
Clashscore	141614	6793 (1.80-1.80)
Ramachandran outliers	138981	6697 (1.80-1.80)
Sidechain outliers	138945	6696 (1.80-1.80)
RSRZ outliers	127900	5850 (1.80-1.80)

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 <=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	А	403	2% 92%	6% •
1	В	403	2% <b>90</b> %	6% ••
1	С	403	% 93%	5%•
1	D	403	2% 90%	8% •
1	Е	403	% • 93%	



Mol	Chain	Length	Quality of chain	
1	F	403	% 91%	6% •
1	G	403	2% 92%	5% •
1	Н	403	% 93%	



## 2 Entry composition (i)

There are 5 unique types of molecules in this entry. The entry contains 28439 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.

Mol	Chain	Residues		At	oms			ZeroOcc	AltConf	Trace
1	1 1	202	Total	С	Ν	0	S	0	0	0
	A		3124	2014	523	582	5	0	9	0
1	р	202	Total	С	Ν	0	S	0	11	0
	D		3141	2027	528	581	5	0	11	0
1	С	307	Total	С	Ν	0	S	0	19	0
1		591	3181	2054	538	584	5	0	12	0
1	П	207	Total	С	Ν	0	S	0	11	0
1	D	091	3170	2044	535	586	5		11	0
1	F	303	Total	С	Ν	0	S	0	12	0
1	Ľ	090	3150	2032	530	583	5	0		
1	Б	202	Total	С	Ν	0	S	0	10	0
	Г		3135	2022	527	581	5	0	10	0
1	C	202	Total	С	Ν	0	S	0	19	0
		3150	2034	525	586	5	0	12	0	
1	ц	202	Total	С	Ν	0	S	0	10	0
	п		3132	2023	525	579	5		10	0

• Molecule 1 is a protein called Putative mandelate racemase.

There are 88 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
А	-1	MET	-	expression tag	UNP Q5NN22
А	0	SER	-	expression tag	UNP Q5NN22
А	1	LEU	-	expression tag	UNP Q5NN22
А	394	GLU	-	expression tag	UNP Q5NN22
А	395	GLY	-	expression tag	UNP Q5NN22
А	396	HIS	-	expression tag	UNP Q5NN22
А	397	HIS	-	expression tag	UNP Q5NN22
А	398	HIS	-	expression tag	UNP Q5NN22
А	399	HIS	-	expression tag	UNP Q5NN22
А	400	HIS	-	expression tag	UNP Q5NN22
А	401	HIS	-	expression tag	UNP Q5NN22
В	-1	MET	-	expression tag	UNP Q5NN22
В	0	SER	-	expression tag	UNP Q5NN22



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Chain	Residue	Modelled	Actual	Comment	Reference
B	1	LEU	-	expression tag	UNP Q5NN22
B	394	GLU	-	expression tag	UNP Q5NN22
B	395	GLY	-	expression tag	UNP Q5NN22
В	396	HIS	-	expression tag	UNP Q5NN22
В	397	HIS	-	expression tag	UNP Q5NN22
В	398	HIS	-	expression tag	UNP Q5NN22
В	399	HIS	-	expression tag	UNP Q5NN22
В	400	HIS	-	expression tag	UNP Q5NN22
В	401	HIS	-	expression tag	UNP Q5NN22
С	-1	MET	-	expression tag	UNP Q5NN22
С	0	SER	-	expression tag	UNP Q5NN22
С	1	LEU	-	expression tag	UNP Q5NN22
С	394	GLU	-	expression tag	UNP Q5NN22
С	395	GLY	-	expression tag	UNP Q5NN22
С	396	HIS	-	expression tag	UNP Q5NN22
С	397	HIS	-	expression tag	UNP Q5NN22
С	398	HIS	-	expression tag	UNP Q5NN22
С	399	HIS	-	expression tag	UNP Q5NN22
С	400	HIS	-	expression tag	UNP Q5NN22
С	401	HIS	-	expression tag	UNP Q5NN22
D	-1	MET	-	expression tag	UNP Q5NN22
D	0	SER	-	expression tag	UNP Q5NN22
D	1	LEU	-	expression tag	UNP Q5NN22
D	394	GLU	-	expression tag	UNP Q5NN22
D	395	GLY	-	expression tag	UNP Q5NN22
D	396	HIS	-	expression tag	UNP Q5NN22
D	397	HIS	-	expression tag	UNP Q5NN22
D	398	HIS	-	expression tag	UNP Q5NN22
D	399	HIS	-	expression tag	UNP Q5NN22
D	400	HIS	-	expression tag	UNP Q5NN22
D	401	HIS	-	expression tag	UNP Q5NN22
Е	-1	MET	-	expression tag	UNP Q5NN22
Е	0	SER	-	expression tag	UNP Q5NN22
Е	1	LEU	-	expression tag	UNP Q5NN22
Е	394	GLU	-	expression tag	UNP Q5NN22
Е	395	GLY	-	expression tag	UNP Q5NN22
Е	396	HIS	-	expression tag	UNP Q5NN22
Е	397	HIS	-	expression tag	UNP Q5NN22
Е	398	HIS	-	expression tag	UNP Q5NN22
Е	399	HIS	-	expression tag	UNP Q5NN22
Е	400	HIS	-	expression tag	UNP Q5NN22
Е	401	HIS	-	expression tag	UNP Q5NN22



Chain	Residue	Modelled	Actual	Comment	Reference
F	-1	MET	-	expression tag	UNP Q5NN22
F	0	SER	-	expression tag	UNP Q5NN22
F	1	LEU	-	expression tag	UNP Q5NN22
F	394	GLU	-	expression tag	UNP Q5NN22
F	395	GLY	-	expression tag	UNP Q5NN22
F	396	HIS	-	expression tag	UNP Q5NN22
F	397	HIS	_	expression tag	UNP Q5NN22
F	398	HIS	-	expression tag	UNP Q5NN22
F	399	HIS	-	expression tag	UNP Q5NN22
F	400	HIS	-	expression tag	UNP Q5NN22
F	401	HIS	-	expression tag	UNP Q5NN22
G	-1	MET	-	expression tag	UNP Q5NN22
G	0	SER	-	expression tag	UNP Q5NN22
G	1	LEU	-	expression tag	UNP Q5NN22
G	394	GLU	-	expression tag	UNP Q5NN22
G	395	GLY	-	expression tag	UNP Q5NN22
G	396	HIS	-	expression tag	UNP Q5NN22
G	397	HIS	-	expression tag	UNP Q5NN22
G	398	HIS	-	expression tag	UNP Q5NN22
G	399	HIS	-	expression tag	UNP Q5NN22
G	400	HIS	-	expression tag	UNP Q5NN22
G	401	HIS	-	expression tag	UNP Q5NN22
Н	-1	MET	-	expression tag	UNP Q5NN22
Н	0	SER	-	expression tag	UNP Q5NN22
Н	1	LEU	-	expression tag	UNP Q5NN22
Н	394	GLU	-	expression tag	UNP Q5NN22
Н	395	GLY	-	expression tag	UNP Q5NN22
Н	396	HIS	-	expression tag	UNP Q5NN22
H	397	HIS	-	expression tag	UNP Q5NN22
H	398	HIS	-	expression tag	UNP Q5NN22
Н	399	HIS	-	expression tag	UNP $Q5NN22$
H	400	HIS	-	expression tag	UNP Q5NN22
H	401	HIS	-	expression tag	UNP Q5NN22

• Molecule 2 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	А	1	Total Mg 1 1	0	0
2	В	1	Total Mg 1 1	0	0
2	С	1	Total Mg 1 1	0	0

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Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	D	1	Total Mg 1 1	0	0
2	Е	1	Total Mg 1 1	0	0
2	F	1	Total Mg 1 1	0	0
2	G	1	Total Mg 1 1	0	0
2	Н	1	Total Mg 1 1	0	0

• Molecule 3 is GLYCEROL (three-letter code: GOL) (formula:  $C_3H_8O_3$ ).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	А	1	$\begin{array}{ccc} \text{Total} & \text{C} & \text{O} \\ 6 & 3 & 3 \end{array}$	0	0
3	А	1	$\begin{array}{ccc} \text{Total} & \text{C} & \text{O} \\ 6 & 3 & 3 \end{array}$	0	0
3	В	1	$\begin{array}{ccc} \text{Total}  \text{C}  \text{O} \\ 6  3  3 \end{array}$	0	0
3	С	1	$\begin{array}{ccc} \text{Total}  \text{C}  \text{O} \\ 6  3  3 \end{array}$	0	0
3	С	1	$\begin{array}{ccc} \text{Total}  \text{C}  \text{O} \\ 6  3  3 \end{array}$	0	0
3	D	1	$\begin{array}{ccc} \text{Total} & \text{C} & \text{O} \\ 6 & 3 & 3 \end{array}$	0	0



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Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	D	1	$\begin{array}{ccc} \text{Total} & \text{C} & \text{O} \\ 6 & 3 & 3 \end{array}$	0	0
3	Е	1	$\begin{array}{ccc} \text{Total}  \text{C}  \text{O} \\ 6  3  3 \end{array}$	0	0
3	F	1	$\begin{array}{ccc} \text{Total}  \text{C}  \text{O} \\ 6  3  3 \end{array}$	0	0
3	F	1	$\begin{array}{ccc} \text{Total}  \text{C}  \text{O} \\ 6  3  3 \end{array}$	0	0
3	G	1	$\begin{array}{ccc} \text{Total}  \text{C}  \text{O} \\ 6  3  3 \end{array}$	0	0
3	G	1	$\begin{array}{ccc} \text{Total}  \text{C}  \text{O} \\ 6  3  3 \end{array}$	0	0
3	G	1	$\begin{array}{ccc} \text{Total} & \text{C} & \text{O} \\ 6 & 3 & 3 \end{array}$	0	0
3	Н	1	$\begin{array}{ccc} \text{Total}  \text{C}  \text{O} \\ 6  3  3 \end{array}$	0	0

• Molecule 4 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	В	1	Total Cl 1 1	0	0
4	D	1	Total Cl 1 1	0	0
4	F	1	Total Cl 1 1	0	0
4	G	1	Total Cl 1 1	0	0

• Molecule 5 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	А	384	Total O 384 384	0	0
5	В	400	Total         O           400         400	0	0
5	С	387	Total O 387 387	0	0
5	D	397	Total O 397 397	0	0
5	Е	413	Total         O           413         413	0	0



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Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	F	394	Total O 394 394	0	0
5	G	392	Total O 392 392	0	0
5	Н	393	Total O 393 393	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: Putative mandelate racemase

• Molecule 1: Putative mandelate racemase









## 4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants	190.18Å $190.41$ Å $85.92$ Å	Deperitor
a, b, c, $\alpha$ , $\beta$ , $\gamma$	$90.00^{\circ}$ $90.00^{\circ}$ $90.00^{\circ}$	Depositor
$\mathbf{P}_{\text{assolution}}(\hat{\mathbf{A}})$	20.00 - 1.80	Depositor
Resolution (A)	34.21 - 1.80	EDS
% Data completeness	99.6 (20.00-1.80)	Depositor
(in resolution range)	99.6 (34.21-1.80)	EDS
R <sub>merge</sub>	0.09	Depositor
R <sub>sym</sub>	0.07	Depositor
$< I/\sigma(I) > 1$	$1.73 (at 1.79 \text{\AA})$	Xtriage
Refinement program	REFMAC 5.3.0028	Depositor
D D.	0.144 , $0.186$	Depositor
$\Pi, \Pi_{free}$	0.144 , $0.185$	DCC
$R_{free}$ test set	8765 reflections $(3.04%)$	wwPDB-VP
Wilson B-factor $(Å^2)$	15.6	Xtriage
Anisotropy	0.104	Xtriage
Bulk solvent $k_{sol}(e/A^3), B_{sol}(A^2)$	0.37, $61.1$	EDS
L-test for twinning <sup>2</sup>	$< L >=0.50, < L^2>=0.33$	Xtriage
Estimated twinning fraction	0.003 for k,h,-l	Xtriage
$F_o, F_c$ correlation	0.97	EDS
Total number of atoms	28439	wwPDB-VP
Average B, all atoms $(Å^2)$	18.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: The largest off-origin peak in the Patterson function is 2.97% of the height of the origin peak. No significant pseudotranslation is detected.

<sup>&</sup>lt;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.



<sup>&</sup>lt;sup>1</sup>Intensities estimated from amplitudes.

# 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, MG, CL

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).

Mal	Chain	Bond lengths		Bond angles	
MIOI	Unain	RMSZ	# Z  > 5	RMSZ	# Z  > 5
1	А	0.47	0/3219	0.58	0/4354
1	В	0.50	0/3242	0.59	0/4382
1	С	0.47	0/3284	0.57	0/4436
1	D	0.49	0/3273	0.58	0/4423
1	Е	0.52	0/3255	0.59	0/4401
1	F	0.49	0/3233	0.58	0/4371
1	G	0.50	0/3245	0.59	1/4389~(0.0%)
1	Н	0.48	0/3230	0.57	0/4367
All	All	0.49	0/25981	0.58	1/35123~(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
1	А	0	1
1	В	0	3
1	С	0	1
1	D	0	1
1	F	0	2
1	G	0	2
1	Н	0	1
All	All	0	11

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	$Observed(^{o})$	$Ideal(^{o})$
1	G	138	GLU	N-CA-C	-5.03	97.43	111.00



There are no chirality outliers.

5 of 11 planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	А	135	TRP	Peptide
1	В	137	LYS	Peptide
1	В	346	LEU	Peptide
1	В	391	GLU	Peptide
1	С	135	TRP	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	А	3124	0	3137	24	0
1	В	3141	0	3169	39	0
1	С	3181	0	3209	24	0
1	D	3170	0	3184	23	0
1	Е	3150	0	3170	23	0
1	F	3135	0	3156	26	0
1	G	3150	0	3170	20	0
1	Н	3132	0	3161	22	0
2	А	1	0	0	0	0
2	В	1	0	0	0	0
2	С	1	0	0	0	0
2	D	1	0	0	0	0
2	Е	1	0	0	0	0
2	F	1	0	0	0	0
2	G	1	0	0	0	0
2	Н	1	0	0	0	0
3	А	12	0	16	0	0
3	В	6	0	8	0	0
3	С	12	0	16	1	0
3	D	12	0	16	1	0
3	Е	6	0	8	0	0
3	F	12	0	16	0	0
3	G	18	0	24	2	0
3	Н	6	0	8	0	0
4	В	1	0	0	0	0
4	D	1	0	0	0	0



Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	F	1	0	0	0	0
4	G	1	0	0	0	0
5	А	384	0	0	0	0
5	В	400	0	0	1	0
5	С	387	0	0	1	0
5	D	397	0	0	1	0
5	Ε	413	0	0	2	0
5	F	394	0	0	3	0
5	G	392	0	0	0	0
5	Н	393	0	0	0	0
All	All	28439	0	25468	142	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 3.

The worst 5 of 142 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:304[B]:MET:CE	1:C:304[B]:MET:HG3	1.82	1.09
1:B:307[B]:ILE:CD1	1:C:304[B]:MET:HE1	1.93	0.97
1:A:307[B]:ILE:CD1	1:H:304[B]:MET:HE1	1.96	0.95
1:C:299:LYS:HE2	3:C:3403:GOL:H31	1.48	0.95
1:B:304[B]:MET:HE1	1:C:304[B]:MET:HG3	1.46	0.94

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	Perce	ntiles
1	А	399/403~(99%)	379~(95%)	19 (5%)	1 (0%)	41	27
1	В	402/403~(100%)	381~(95%)	19 (5%)	2 (0%)	29	15



Mol	Chain	Analysed	Favoured	Allowed	Outliers	Perce	ntiles
1	С	405/403~(100%)	386~(95%)	18 (4%)	1 (0%)	47	33
1	D	404/403~(100%)	390~(96%)	13 (3%)	1 (0%)	47	33
1	Е	403/403~(100%)	385~(96%)	17 (4%)	1 (0%)	47	33
1	F	401/403~(100%)	384 (96%)	16 (4%)	1 (0%)	47	33
1	G	403/403 (100%)	386~(96%)	16 (4%)	1 (0%)	47	33
1	Н	401/403~(100%)	384 (96%)	15 (4%)	2(0%)	29	15
All	All	3218/3224 (100%)	3075~(96%)	133 (4%)	10 (0%)	41	27

5 of 10 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	В	344	LYS
1	С	136	GLY
1	D	264	GLY
1	Е	264	GLY
1	Н	136	GLY

#### 5.3.2 Protein sidechains (i)

In the following table, the Percentiles column shows the percent side chain 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	А	329/329~(100%)	323~(98%)	6(2%)	59 48
1	В	331/329~(101%)	327~(99%)	4 (1%)	71 65
1	С	335/329~(102%)	326~(97%)	9~(3%)	44 31
1	D	334/329~(102%)	326~(98%)	8 (2%)	49 36
1	Ε	332/329~(101%)	325~(98%)	7(2%)	53 42
1	F	330/329~(100%)	323~(98%)	7~(2%)	53 42
1	G	332/329~(101%)	325~(98%)	7(2%)	53 42
1	Н	330/329~(100%)	323~(98%)	7(2%)	53 42
All	All	2653/2632~(101%)	2598~(98%)	55 (2%)	53 42



5 of 55 residues with a non-rotameric sidechain are listed below:

Mol	Chain	$\mathbf{Res}$	Type
1	Ε	131	LEU
1	F	137	LYS
1	Н	391	GLU
1	Н	137	LYS
1	Е	137	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 28 such sidechains are listed below:

Mol	Chain	$\mathbf{Res}$	Type
1	D	398	HIS
1	Н	343	GLN
1	Е	234	ASN
1	Н	173	ASN
1	Е	214	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)

Of 26 ligands modelled in this entry, 12 are monoatomic - leaving 14 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).



Mal	Type	Chain	Dog	Link	B	ond leng	$_{ m gths}$	В	Sond ang	gles
WIOI	туре	Ullalli	nes		Counts	RMSZ	# Z >2	Counts	RMSZ	# Z >2
3	GOL	А	3414	-	$5,\!5,\!5$	0.32	0	$5,\!5,\!5$	0.59	0
3	GOL	D	3406	-	$5,\!5,\!5$	0.35	0	$5,\!5,\!5$	0.50	0
3	GOL	А	3401	-	$5,\!5,\!5$	0.33	0	$5,\!5,\!5$	0.40	0
3	GOL	В	3402	-	$5,\!5,\!5$	0.33	0	$5,\!5,\!5$	0.48	0
3	GOL	F	3409	-	$5,\!5,\!5$	0.35	0	$5,\!5,\!5$	0.55	0
3	GOL	G	3405	-	$5,\!5,\!5$	0.35	0	$5,\!5,\!5$	0.58	0
3	GOL	Н	3410	-	$5,\!5,\!5$	0.34	0	$5,\!5,\!5$	0.46	0
3	GOL	G	3408	-	$5,\!5,\!5$	0.35	0	$5,\!5,\!5$	0.38	0
3	GOL	Е	3407	-	$5,\!5,\!5$	0.38	0	$5,\!5,\!5$	0.48	0
3	GOL	G	3412	-	$5,\!5,\!5$	0.41	0	$5,\!5,\!5$	0.66	0
3	GOL	С	3403	-	$5,\!5,\!5$	0.39	0	$5,\!5,\!5$	0.31	0
3	GOL	D	3413	-	$5,\!5,\!5$	0.33	0	$5,\!5,\!5$	0.69	0
3	GOL	F	3411	-	$5,\!5,\!5$	0.39	0	$5,\!5,\!5$	0.48	0
3	GOL	С	3404	-	$5,\!5,\!5$	0.44	0	$5,\!5,\!5$	0.44	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	GOL	А	3414	-	-	2/4/4/4	-
3	GOL	D	3406	-	-	0/4/4/4	-
3	GOL	А	3401	-	-	0/4/4/4	-
3	GOL	В	3402	-	-	0/4/4/4	-
3	GOL	F	3409	-	-	2/4/4/4	-
3	GOL	G	3405	-	-	0/4/4/4	-
3	GOL	Н	3410	-	-	0/4/4/4	-
3	GOL	G	3408	-	-	4/4/4/4	-
3	GOL	Е	3407	-	-	2/4/4/4	-
3	GOL	G	3412	-	-	4/4/4/4	-
3	GOL	С	3403	-	-	2/4/4/4	-
3	GOL	D	3413	-	-	0/4/4/4	-
3	GOL	F	3411	-	-	4/4/4/4	-
3	GOL	C	3404	-	-	0/4/4/4	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

5 of 20 torsion outliers are listed below:



Mol	Chain	Res	Type	Atoms
3	А	3414	GOL	C1-C2-C3-O3
3	С	3403	GOL	C1-C2-C3-O3
3	Е	3407	GOL	O1-C1-C2-C3
3	F	3409	GOL	O1-C1-C2-C3
3	F	3411	GOL	C1-C2-C3-O3

There are no ring outliers.

3 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	G	3412	GOL	2	0
3	С	3403	GOL	1	0
3	D	3413	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 (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,  $95^{th}$  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	А	393/403~(97%)	-0.44	8 (2%) 65 61	8, 15, 38, 71	0
1	В	393/403~(97%)	-0.49	10 (2%) 57 52	7, 14, 40, 70	0
1	С	397/403~(98%)	-0.42	6 (1%) 73 70	8, 14, 35, 74	0
1	D	397/403~(98%)	-0.46	8 (2%) 65 61	7, 14, 37, 73	0
1	Ε	393/403~(97%)	-0.55	6 (1%) 73 70	6, 12, 32, 68	0
1	F	393/403~(97%)	-0.54	4 (1%) 82 80	7, 14, 32, 58	0
1	G	393/403~(97%)	-0.53	8 (2%) 65 61	7, 13, 40, 65	0
1	Н	393/403~(97%)	-0.52	6 (1%) 73 70	7, 13, 32, 65	0
All	All	3152/3224 (97%)	-0.49	56 (1%) 68 64	6, 14, 36, 74	0

The worst 5 of 56 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	В	346	LEU	6.5
1	А	346	LEU	6.3
1	В	392	ALA	5.8
1	В	347	LEU	5.1
1	D	397	HIS	5.0

## 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,  $95^{th}$  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	$B-factors(A^2)$	Q<0.9
3	GOL	А	3414	6/6	0.82	0.21	27,45,48,50	0
3	GOL	D	3413	6/6	0.85	0.20	37,38,50,55	0
3	GOL	G	3408	6/6	0.87	0.20	29,38,41,49	0
3	GOL	С	3403	6/6	0.89	0.19	29,33,60,62	0
3	GOL	F	3411	6/6	0.91	0.16	22,35,46,51	0
3	GOL	Е	3407	6/6	0.93	0.14	21,25,28,30	0
3	GOL	Н	3410	6/6	0.93	0.13	19,24,26,31	0
3	GOL	G	3405	6/6	0.94	0.10	20,25,33,35	0
3	GOL	А	3401	6/6	0.94	0.09	16,18,24,24	0
3	GOL	В	3402	6/6	0.94	0.10	21,29,31,34	0
3	GOL	G	3412	6/6	0.95	0.16	23,30,34,42	0
3	GOL	D	3406	6/6	0.95	0.09	14,20,25,25	0
3	GOL	F	3409	6/6	0.96	0.08	18,22,25,33	0
3	GOL	С	3404	6/6	0.96	0.09	14,19,25,25	0
2	MG	D	402	1/1	0.99	0.07	8,8,8,8	0
4	CL	В	501	1/1	0.99	0.06	12,12,12,12	0
4	CL	G	501	1/1	0.99	0.04	12,12,12,12	0
2	MG	А	402	1/1	1.00	0.07	10,10,10,10	0
2	MG	Е	402	1/1	1.00	0.10	6,6,6,6	0
2	MG	F	402	1/1	1.00	0.09	8,8,8,8	0
2	MG	G	502	1/1	1.00	0.11	8,8,8,8	0
2	MG	Н	402	1/1	1.00	0.11	8,8,8,8	0
2	MG	В	502	1/1	1.00	0.06	9,9,9,9	0
4	CL	D	501	1/1	1.00	0.06	12,12,12,12	0
4	CL	F	501	1/1	1.00	0.04	11,11,11,11	0
2	MG	С	402	1/1	1.00	0.11	8,8,8,8	0

### 6.5 Other polymers (i)

There are no such residues in this entry.

