wwPDB X-ray Structure Validation Summary Report (i)

May 14, 2020 – 12:02 pm BST

PDB ID	:	5GUQ
Title	:	Crystal structure of ASCH from Zymomonas mobilis
Authors	:	Ha, S.C.; Park, S.Y.; Kim, J.S.
Deposited on	:	2016-08-30
Resolution	:	1.70 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 following versions of software and data (see references (1)) were used in the production of this report:

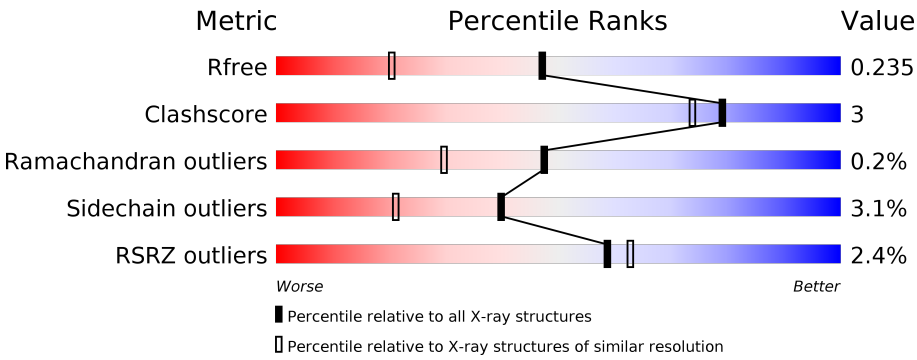
MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.11
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.11

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 1.70 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	4298 (1.70-1.70)
Clashscore	141614	4695 (1.70-1.70)
Ramachandran outliers	138981	4610 (1.70-1.70)
Sidechain outliers	138945	4610 (1.70-1.70)
RSRZ outliers	127900	4222 (1.70-1.70)

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 >=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	A	149	3% 86% 9% • •
1	B	149	2% 87% 7% • •
1	C	149	2% 93% 6% •
1	D	149	2% 88% 7% • •

2 Entry composition (i)

There are 2 unique types of molecules in this entry. The entry contains 5566 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 Helix-turn-helix domain-containing protein.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf	Trace
1	A	143	Total C N O 1187 764 210 213	0	1	0
1	B	143	Total C N O 1200 773 211 216	0	3	0
1	C	149	Total C N O S 1253 807 221 224 1	0	4	0
1	D	143	Total C N O 1196 772 210 214	0	3	0

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	0	GLY	-	expression tag	UNP A0A0H3G0N3
B	0	GLY	-	expression tag	UNP A0A0H3G0N3
C	0	GLY	-	expression tag	UNP A0A0H3G0N3
D	0	GLY	-	expression tag	UNP A0A0H3G0N3

- Molecule 2 is water.

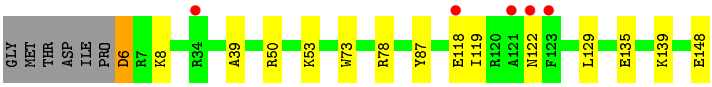
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	158	Total O 158 158	0	0
2	B	163	Total O 163 163	0	0
2	C	235	Total O 235 235	0	0
2	D	174	Total O 174 174	0	0

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.

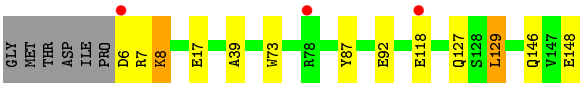
- Molecule 1: Helix-turn-helix domain-containing protein

Chain A:



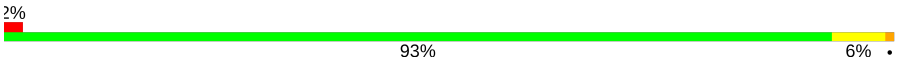
- Molecule 1: Helix-turn-helix domain-containing protein

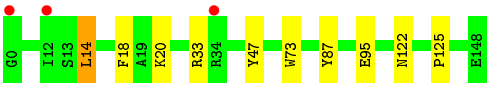
Chain B:



- Molecule 1: Helix-turn-helix domain-containing protein

Chain C:





- Molecule 1: Helix-turn-helix domain-containing protein

Chain D:

4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α, β, γ	72.91Å 52.86Å 88.67Å 90.00° 111.56° 90.00°	Depositor
Resolution (Å)	30.59 – 1.70 30.59 – 1.70	Depositor EDS
% Data completeness (in resolution range)	98.7 (30.59-1.70) 98.7 (30.59-1.70)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	2.71 (at 1.70Å)	Xtriage
Refinement program	PHENIX 1.8.1_1168	Depositor
R, R_{free}	0.198 , 0.235 0.199 , 0.235	Depositor DCC
R_{free} test set	2751 reflections (4.00%)	wwPDB-VP
Wilson B-factor (Å^2)	21.0	Xtriage
Anisotropy	0.401	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.34 , 48.7	EDS
L-test for twinning[2]	$< \|L\| > = 0.48$, $< L^2 > = 0.31$	Xtriage
Estimated twinning fraction	0.023 for h,-k,-h-l	Xtriage
F_o,F_c correlation	0.95	EDS
Total number of atoms	5566	wwPDB-VP
Average B, all atoms (Å^2)	25.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 35.02 % 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 6.2234e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ 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)

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	0.37	0/1215	0.59	0/1646
1	B	0.35	0/1231	0.60	1/1668 (0.1%)
1	C	0.40	0/1285	0.64	0/1741
1	D	0.37	0/1230	0.59	0/1667
All	All	0.37	0/4961	0.60	1/6722 (0.0%)

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	B	129	LEU	CA-CB-CG	6.19	129.53	115.30

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	1187	0	1203	7	0
1	B	1200	0	1217	8	0
1	C	1253	0	1280	5	0
1	D	1196	0	1221	7	0
2	A	158	0	0	2	0
2	B	163	0	0	3	0
2	C	235	0	0	1	0
2	D	174	0	0	1	1

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
All	All	5566	0	4921	27	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 3.

The worst 5 of 27 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:119:ILE:HG23	1:A:139:LYS:HE3	1.67	0.75
1:A:6:ASP:N	2:A:201:HOH:O	2.25	0.68
1:C:122:ASN:ND2	2:C:203:HOH:O	2.27	0.68
1:A:78:ARG:NH1	2:A:204:HOH:O	2.38	0.57
1:D:53:LYS:HE2	1:D:113:LEU:HD12	1.88	0.56

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)
2:D:360:HOH:O	2:D:369:HOH:O[2_444]	2.17	0.03

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	142/149 (95%)	137 (96%)	5 (4%)	0	100 100
1	B	144/149 (97%)	141 (98%)	3 (2%)	0	100 100
1	C	151/149 (101%)	147 (97%)	3 (2%)	1 (1%)	22 8
1	D	144/149 (97%)	141 (98%)	3 (2%)	0	100 100
All	All	581/596 (98%)	566 (97%)	14 (2%)	1 (0%)	47 30

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	33	ARG

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	127/131 (97%)	122 (96%)	5 (4%)	32	13
1	B	129/131 (98%)	124 (96%)	5 (4%)	32	13
1	C	135/131 (103%)	132 (98%)	3 (2%)	52	34
1	D	129/131 (98%)	125 (97%)	4 (3%)	40	21
All	All	520/524 (99%)	503 (97%)	17 (3%)	40	19

5 of 17 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	118	GLU
1	B	129	LEU
1	D	17	GLU
1	B	17	GLU
1	D	34	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

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)

There are no ligands in this entry.

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, 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	143/149 (95%)	0.02	5 (3%)	44	49	16, 23, 45, 66	0
1	B	143/149 (95%)	0.01	3 (2%)	63	67	16, 23, 41, 75	0
1	C	149/149 (100%)	0.01	3 (2%)	65	69	12, 16, 31, 56	0
1	D	143/149 (95%)	0.04	3 (2%)	63	67	14, 21, 40, 54	0
All	All	578/596 (96%)	0.02	14 (2%)	59	63	12, 21, 39, 75	0

The worst 5 of 14 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	6	ASP	6.2
1	C	0	GLY	5.8
1	A	122	ASN	3.9
1	D	6	ASP	3.6
1	C	34	ARG	3.5

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)

There are no ligands in this entry.

6.5 Other polymers (i)

There are no such residues in this entry.