



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 12, 2017 – 10:26 pm GMT

PDB ID : 3EY7
Title : Structure from the mobile metagenome of *V. Cholerae*. Integron cassette protein VCH_CASS1
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Deposited on : 2008-10-20
Resolution : 1.60 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

<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:

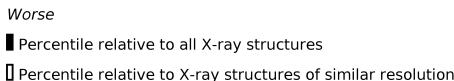
MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.9-1692
EDS	:	trunk28620
Percentile statistics	:	20161228.v01 (using entries in the PDB archive December 28th 2016)
Refmac	:	5.8.0135
CCP4	:	6.5.0
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	recalc28949

i

X-RAY DIFFRACTION

A.

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	2696 (1.60-1.60)
Clashscore	112137	2967 (1.60-1.60)
Ramachandran outliers	110173	2887 (1.60-1.60)
Sidechain outliers	110143	2886 (1.60-1.60)
RSRZ outliers	101464	2714 (1.60-1.60)

electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	133	<div> <div></div> <div>92%</div> <div>7%</div> </div>
1	B	133	<div> <div></div> <div>85%</div> <div>11%</div> </div>

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 2515 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 Biphenyl-2,3-diol 1,2-dioxygenase III-related protein.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	132	Total	C	N	O	S	Se	0	3	0
			1043	661	176	199	1	6			
1	B	128	Total	C	N	O	S	Se	0	6	0
			1038	662	175	195	1	5			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	0	GLY	-	EXPRESSION TAG	UNP Q9K3D3
A	92	ASP	ASN	ENGINEERED	UNP Q9K3D3
B	0	GLY	-	EXPRESSION TAG	UNP Q9K3D3
B	92	ASP	ASN	ENGINEERED	UNP Q9K3D3

- Molecule 2 is CALCIUM ION (three-letter code: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	1	Total	Ca	0	0
			1	1		
2	A	1	Total	Ca	0	0
			1	1		

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	221	Total	O	0	0
			221	221		
3	B	211	Total	O	0	0
			211	211		

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: Biphenyl-2,3-diol 1,2-dioxygenase III-related protein

Chain A: 



- Molecule 1: Biphenyl-2,3-diol 1,2-dioxygenase III-related protein

Chain B: 



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	43.03Å 67.87Å 91.24Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	27.75 – 1.60 27.75 – 1.60	Depositor EDS
% Data completeness (in resolution range)	96.3 (27.75-1.60) 96.3 (27.75-1.60)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	0.08	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.06 (at 1.60Å)	Xtriage
Refinement program	PHENIX (phenix.refine)	Depositor
R, R_{free}	0.171 , 0.216 0.168 , 0.215	Depositor DCC
R_{free} test set	1998 reflections (5.77%)	DCC
Wilson B-factor (Å ²)	16.8	Xtriage
Anisotropy	0.556	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 47.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	2515	wwPDB-VP
Average B, all atoms (Å ²)	22.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.41% 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: CA

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	0.35	0/1057	0.54	0/1419
1	B	0.37	0/1054	0.55	0/1419
All	All	0.36	0/2111	0.54	0/2838

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	1043	0	1032	7	0
1	B	1038	0	1030	12	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	221	0	0	2	0
3	B	211	0	0	0	0
All	All	2515	0	2062	18	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.

All (18) 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:43:ILE:HD11	3:A:298:HOH:O	1.96	0.64
1:A:25:ASN:HD21	1:A:29:LYS:HE3	1.67	0.59
1:B:87[B]:MSE:HG2	1:B:114:PHE:CD1	2.40	0.56
1:A:20:ILE:O	1:A:24:THR:HG23	2.08	0.54
1:B:27:TYR:CD2	1:B:31[A]:LEU:HD12	2.44	0.52
1:A:87[A]:MSE:HG2	1:A:114:PHE:CD2	2.45	0.52
1:A:42:ARG:HD3	3:A:177:HOH:O	2.12	0.50
1:B:51:LYS:NZ	1:B:53:ASN:HD21	2.08	0.50
1:B:55[B]:HIS:HE1	1:B:59:ASN:O	1.94	0.49
1:A:87[B]:MSE:HE2	1:A:114:PHE:CD2	2.49	0.48
1:A:11:ASP:HB2	1:B:78:ILE:HG12	1.97	0.47
1:B:84:SER:HA	1:B:87[B]:MSE:HE3	1.96	0.46
1:B:42:ARG:HE	1:B:61:PHE:HE1	1.64	0.46
1:B:51:LYS:HZ2	1:B:53:ASN:HD21	1.64	0.44
1:B:14:VAL:HG22	1:B:53:ASN:HB2	2.03	0.41
1:B:55[B]:HIS:CE1	1:B:59:ASN:O	2.73	0.41
1:B:87[B]:MSE:HG2	1:B:114:PHE:CG	2.55	0.41
1:B:118:ASP:HB2	1:B:119:PRO:CD	2.51	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	133/133 (100%)	129 (97%)	4 (3%)	0	100	100
1	B	132/133 (99%)	130 (98%)	2 (2%)	0	100	100
All	All	265/266 (100%)	259 (98%)	6 (2%)	0	100	100

There are no Ramachandran outliers to report.

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	114/107 (106%)	114 (100%)	0	100	100
1	B	114/107 (106%)	114 (100%)	0	100	100
All	All	228/214 (106%)	228 (100%)	0	100	100

There are no protein residues with a non-rotameric sidechain to report.

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	25	ASN
1	B	53	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 2 ligands modelled in this entry, 2 are monoatomic - leaving 0 for Mogul analysis.

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.

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 [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	127/133 (95%)	-0.06	1 (0%) 86 87	10, 19, 37, 51	0
1	B	124/133 (93%)	-0.09	5 (4%) 39 37	10, 16, 39, 67	0
All	All	251/266 (94%)	-0.07	6 (2%) 59 59	10, 17, 38, 67	0

All (6) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	61	PHE	4.1
1	B	60	GLU	3.5
1	B	3	PHE	2.7
1	B	40	ALA	2.5
1	A	15[A]	LEU	2.4
1	B	130	SER	2.1

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, 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	LLDF	B-factors(\AA^2)	Q<0.9
2	CA	B	133	1/1	0.99	0.04	-2.71	17,17,17,17	0
2	CA	A	133	1/1	0.99	0.04	-3.28	24,24,24,24	0

6.5 Other polymers [i](#)

There are no such residues in this entry.