

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

#### May 22, 2020 – 12:21 pm BST

PDB ID	:	$5\mathrm{E1W}$
$\operatorname{Title}$	:	Crystal structure of the organohalide sensing RdhR-CbdbA1625 transcrip-
		tional regulator in the ligand free form
Authors	:	Quezada, C.P.; Dunstan, M.S.; Fisher, K.; Leys, D.
Deposited on	:	2015-09-30
Resolution	:	1.39 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 (i) symbol.

The following versions of software and data (see references (1)) were used in the production of this report:

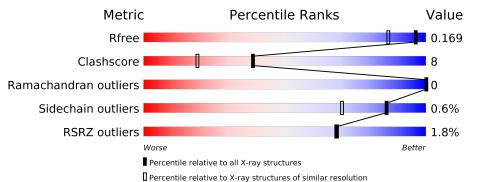
$\operatorname{MolProbity}$	:	4.02b-467
Xtriage (Phenix)	:	1.13
$\mathrm{EDS}$	:	2.11
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
$\operatorname{Refmac}$	:	5.8.0158
$\operatorname{CCP4}$	:	$7.0.044 (\mathrm{Gargrove})$
Ideal geometry (proteins)	:	Engh & 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 1.39 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	$egin{array}{c} { m Whole \ archive} \ (\#{ m Entries}) \end{array}$	${f Similar\ resolution}\ (\#{ m Entries},{ m resolution\ range}({ m \AA}))$
$R_{free}$	130704	1714(1.40-1.40)
Clashscore	141614	1812 (1.40-1.40)
Ramachandran outliers	138981	1763(1.40-1.40)
Sidechain outliers	138945	1762(1.40-1.40)
RSRZ outliers	127900	1674(1.40-1.40)

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			
			2%			
1	A	180	81%	10%	• 7%	-



# 2 Entry composition (i)

There are 2 unique types of molecules in this entry. The entry contains 1723 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 Transcriptional regulator, MarR family.

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace		
1	А	168	Total 1475	m C 951	N 254	O 263	${ m S}$ 7	0	17	0

Chain	Residue	Modelled	Actual	Comment	Reference
A	-18	MET	-	initiating methionine	UNP M1R817
A	-17	ALA	-	expression tag	UNP M1R817
A	-16	HIS	-	expression tag	UNP M1R817
A	-15	HIS	-	expression tag	UNP M1R817
A	-14	HIS	-	expression tag	UNP M1R817
A	-13	HIS	-	expression tag	UNP M1R817
A	-12	HIS	-	expression tag	UNP M1R817
A	-11	HIS	-	expression tag	UNP M1R817
A	-10	SER	-	expression tag	UNP M1R817
A	-9	SER	-	expression tag	UNP M1R817
A	-8	GLY	-	expression tag	UNP M1R817
A	-7	LEU	-	expression tag	UNP M1R817
A	-6	GLU	-	expression tag	UNP M1R817
A	-5	VAL	-	expression tag	UNP M1R817
A	-4	LEU	-	expression tag	UNP M1R817
A	-3	PHE	-	expression tag	UNP M1R817
А	-2	GLN	-	expression tag	UNP M1R817
A	-1	GLY	-	expression tag	UNP M1R817
A	1	PRO	-	expression tag	UNP M1R817
A	105	ILE	VAL	conflict	UNP M1R817

There are 20 discrepancies between the modelled and reference sequences:

• Molecule 2 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	А	248	Total         O           248         248	0	0

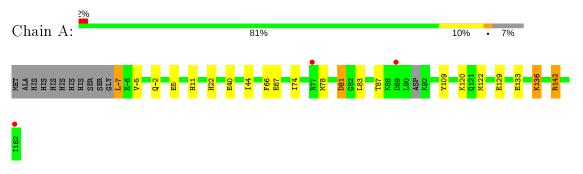




# 3 Residue-property plots (i)

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: Transcriptional regulator, MarR family





## 4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 31 2 1	Depositor
$\begin{array}{c} \text{Cell constants} \\ \text{a, b, c, } \alpha, \beta, \gamma \end{array}$	$\begin{array}{cccc} 71.95 \text{\AA} & 71.95 \text{\AA} & 87.63 \text{\AA} \\ 90.00^\circ & 90.00^\circ & 120.00^\circ \end{array}$	Depositor
Resolution (Å)	$\begin{array}{rrrrrrrrrrrrrrrrrrrrrrrrrrrrrrrrrrrr$	Depositor EDS
% Data completeness (in resolution range)	$\begin{array}{c} 99.8 \ (31.16\text{-}1.39) \\ 99.8 \ (31.16\text{-}1.39) \end{array}$	Depositor EDS
R <sub>merge</sub>	0.05	Depositor
R <sub>sym</sub>	(Not available)	Depositor
$< I/\sigma(I) > 1$	$3.00 ({ m at} 1.39{ m \AA})$	Xtriage
Refinement program	REFMAC $5.8.0049$	Depositor
$R, R_{free}$	$\begin{array}{rrrrrrrrrrrrrrrrrrrrrrrrrrrrrrrrrrrr$	Depositor DCC
$R_{free}$ test set	2688 reflections $(5.08%)$	wwPDB-VP
Wilson B-factor $(Å^2)$	13.2	Xtriage
Anisotropy	0.015	Xtriage
Bulk solvent $k_{sol}(e/Å^3), B_{sol}(Å^2)$	0.35 , 55.7	EDS
L-test for $twinning^2$	$<  L  > = 0.49, < L^2 > = 0.32$	Xtriage
Estimated twinning fraction	0.028 for -h,-k,l	Xtriage
$F_o, F_c$ correlation	0.97	EDS
Total number of atoms	1723	wwPDB-VP
Average B, all atoms $(Å^2)$	19.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: The largest off-origin peak in the Patterson function is 8.42% 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.



 $<sup>^1 {\</sup>rm Intensities}$  estimated from amplitudes.

# 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).

Mal	Chain	Bo	nd lengths	Bond angles		
Mol	Cham	RMSZ	# Z  > 5	RMSZ	# Z  > 5	
1	А	1.37	12/1542~(0.8%)	1.18	8/2072~(0.4%)	

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.

$\mathbf{Mol}$	Chain	#Chirality outliers	#Planarity outliers
1	А	0	1

All $(12)$	$\operatorname{bond}$	length	outliers	$\operatorname{are}$	listed	below:
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Mol	Chain	Res	Type	Atoms		$\operatorname{Observed}(\operatorname{\AA})$	Ideal(Å)
1	А	133[A]	GLU	CD-OE2	16.08	1.43	1.25
1	А	133[B]	GLU	CD-OE2	16.08	1.43	1.25
1	А	40	GLU	CD-OE1	9.55	1.36	1.25
1	А	133[A]	GLU	CG-CD	6.87	1.62	1.51
1	А	133[B]	GLU	CG-CD	6.87	1.62	1.51
1	А	67	GLU	CG-CD	-6.36	1.42	1.51
1	А	129	GLU	CG-CD	5.61	1.60	1.51
1	А	66	PHE	CD1-CE1	5.49	1.50	1.39
1	А	133[A]	GLU	CB-CG	5.30	1.62	1.52
1	А	133[B]	GLU	CB-CG	5.30	1.62	1.52
1	А	5	GLU	CD-OE1	5.28	1.31	1.25
1	А	122	MET	SD-CE	-5.16	1.49	1.77

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	$Observed(^{o})$	$Ideal(^{o})$
1	А	136	LYS	CD-CE-NZ	6.59	126.86	111.70
1	А	122	MET	CG-SD-CE	-6.45	89.88	100.20
1	А	142	ARG	NE-CZ-NH1	-6.41	117.10	120.30
1	А	-7	LEU	CB-CG-CD1	6.38	121.84	111.00

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Mol	Chain	Res	Type	Atoms	Z	$Observed(^{o})$	$Ideal(^{o})$
1	А	-7	LEU	CA-CB-CG	-6.16	101.12	115.30
1	А	122	MET	CB-CG-SD	-5.83	94.90	112.40
1	А	66	PHE	CB-CG-CD1	5.62	124.73	120.80
1	А	81	ASP	CB-CG-OD2	-5.50	113.35	118.30

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There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	А	142	ARG	Sidechain

### 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	1475	0	1568	23	2
2	А	248	0	0	10	6
All	All	1723	0	1568	23	6

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 (23) 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:22[B]:HIS:CD2	2:A:202:HOH:O	1.83	1.25
1:A:22[B]:HIS:HD2	2:A:202:HOH:O	1.28	0.94
1:A:44:ILE:HG21	1:A:78[B]:MET:HE3	1.56	0.87
1:A:-5[A]:VAL:CG1	1:A:-2:GLN:HG3	2.12	0.79
1:A:74:ILE:HG22	1:A:78[B]:MET:HE2	1.66	0.77
1:A:-5[A]:VAL:HG11	1:A:-2:GLN:HG3	1.71	0.72
1:A:87:THR:HG22	2:A:368:HOH:O	1.92	0.68
1:A:74:ILE:HG22	1:A:78[B]:MET:CE	2.25	0.66
1:A:-5[A]:VAL:HG13	1:A:-2:GLN:HG3	1.81	0.62
1:A:87:THR:HG23	2:A:328:HOH:O	2.07	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:44:ILE:HD12	1:A:74:ILE:HG21	1.89	0.53
1:A:136:LYS:HE2	2:A:387:HOH:O	2.08	0.53
1:A:44:ILE:CG2	1:A:78[B]:MET:HE3	2.36	0.50
1:A:11:HIS:HE1	2:A:391:HOH:O	1.93	0.50
1:A:-5[B]:VAL:HG22	2:A:320:HOH:O	2.10	0.50
1:A:-5[B]:VAL:CG2	1:A:-2:GLN:HG3	2.43	0.47
1:A:120[A]:LYS:HG3	2:A:203:HOH:O	2.16	0.45
1:A:-5[A]:VAL:HG13	1:A:-2:GLN:CG	2.46	0.44
1:A:11:HIS:HD2	2:A:339:HOH:O	1.99	0.44
1:A:81:ASP:HB2	1:A:83[B]:LEU:HD23	2.01	0.43
1:A:120[A]:LYS:HE2	2:A:384:HOH:O	2.19	0.43
1:A:44:ILE:HD12	1:A:74:ILE:CG2	2.52	0.40

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All (6) 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	$\begin{array}{c} {\rm Interatomic}\\ {\rm distance}~({\rm \AA}) \end{array}$	Clash overlap (Å)
2:A:302:HOH:O	2:A:395:HOH:O[3_454]	0.33	1.87
2:A:391:HOH:O	2:A:405:HOH:O[6_555]	0.37	1.83
2:A:381:HOH:O	2:A:402:HOH:O[6_555]	0.39	1.81
2:A:345:HOH:O	2:A:419:HOH:O[3_564]	1.96	0.24
1:A:120[A]:LYS:NZ	2:A:228:HOH:O[6_555]	2.09	0.11
1:A:109:TYR:OH	2:A:356:HOH:O[3_454]	2.11	0.09

### 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	А	182/180~(101%)	179~(98%)	3~(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	А	174/167~(104%)	173~(99%)	1 (1%)	86 70

All (1) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	А	-7	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	А	11	HIS

#### 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, 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	$<$ <b>RSRZ</b> $>$	$\# RSRZ {>}2$		$\mathbf{OWAB}(\mathbf{A}^2)$	Q<0.9	
1	А	168/180~(93%)	-0.16	3 (1%)	68	68	10, 14, 32, 52	0

All (3) RSRZ outliers are listed below:

Mol	Chain	$\mathbf{Res}$	Type	RSRZ
1	А	162	ILE	5.5
1	А	77	ARG	2.3
1	А	89	ASP	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)

There are no ligands in this entry.

#### 6.5 Other polymers (i)

There are no such residues in this entry.

