



Full wwPDB X-ray Structure Validation Report ⓘ

May 25, 2020 – 07:08 am BST

PDB ID : 1WLY
Title : Crystal Structure of 2-Haloacrylate Reductase
Authors : Omi, R.
Deposited on : 2004-06-30
Resolution : 1.30 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 ⓘ symbol.

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

MolProbity : 4.02b-467
Xtrriage (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

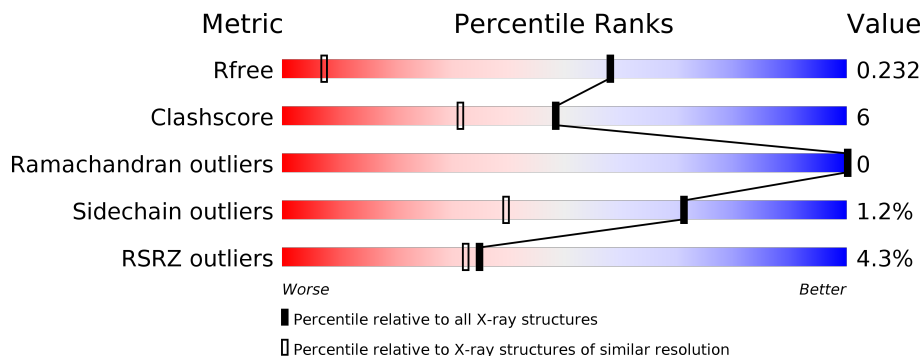
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.30 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	1058 (1.30-1.30)
Clashscore	141614	1101 (1.30-1.30)
Ramachandran outliers	138981	1058 (1.30-1.30)
Sidechain outliers	138945	1058 (1.30-1.30)
RSRZ outliers	127900	1029 (1.30-1.30)

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 $\leq 5\%$. 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	333	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 2765 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 2-haloacrylate reductase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	322	2430	1538	432	446	14	0	0	0

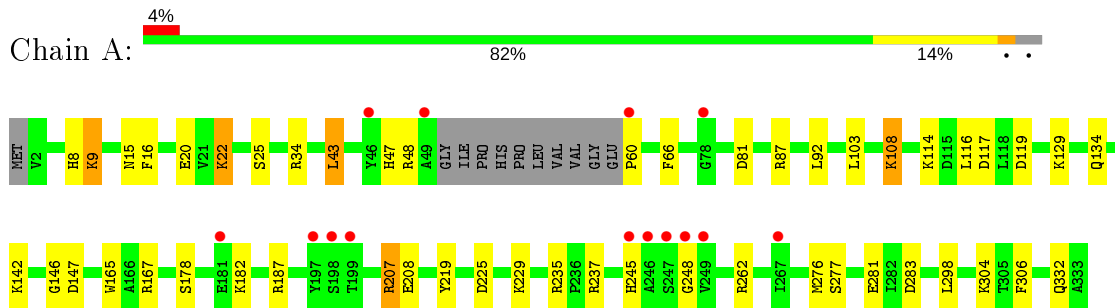
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	335	Total	O	0	0
			335	335		

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: 2-haloacrylate reductase



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	140.80Å 49.18Å 59.05Å 90.00° 100.97° 90.00°	Depositor
Resolution (Å)	19.33 – 1.30 28.75 – 1.30	Depositor EDS
% Data completeness (in resolution range)	97.0 (19.33-1.30) 97.2 (28.75-1.30)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.76 (at 1.30Å)	Xtrriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.212 , 0.232 0.212 , 0.232	Depositor DCC
R_{free} test set	9462 reflections (9.98%)	wwPDB-VP
Wilson B-factor (Å ²)	13.8	Xtrriage
Anisotropy	0.815	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.41 , 47.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	2765	wwPDB-VP
Average B, all atoms (Å ²)	19.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.27% of the height of the origin peak. No significant pseudotranslation is detected.*

¹ Intensities estimated from amplitudes.

² 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

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	1.29	7/2482 (0.3%)	1.29	25/3365 (0.7%)

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	16	PHE	CE2-CZ	6.41	1.49	1.37
1	A	25	SER	CB-OG	-6.05	1.34	1.42
1	A	281	GLU	CD-OE2	6.01	1.32	1.25
1	A	208	GLU	CG-CD	6.00	1.60	1.51
1	A	66	PHE	CD2-CE2	5.15	1.49	1.39
1	A	165	TRP	CE3-CZ3	5.12	1.47	1.38
1	A	219	TYR	CE2-CZ	5.11	1.45	1.38

All (25) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	235	ARG	NE-CZ-NH2	10.47	125.53	120.30
1	A	147	ASP	CB-CG-OD2	-9.07	110.13	118.30
1	A	237	ARG	NE-CZ-NH2	8.50	124.55	120.30
1	A	119	ASP	CB-CG-OD1	7.51	125.06	118.30
1	A	187	ARG	NE-CZ-NH2	-7.42	116.59	120.30
1	A	237	ARG	NE-CZ-NH1	-7.20	116.70	120.30
1	A	87	ARG	NE-CZ-NH2	-7.15	116.72	120.30
1	A	262	ARG	NE-CZ-NH2	7.08	123.84	120.30
1	A	167	ARG	NE-CZ-NH2	-6.86	116.87	120.30
1	A	108	LYS	CD-CE-NZ	6.86	127.47	111.70
1	A	225	ASP	CB-CG-OD2	-6.35	112.58	118.30
1	A	298	LEU	CB-CA-C	-6.26	98.31	110.20
1	A	208	GLU	OE1-CD-OE2	-6.09	116.00	123.30
1	A	34	ARG	NE-CZ-NH1	-6.00	117.30	120.30
1	A	117	ASP	CB-CG-OD2	-5.90	112.99	118.30
1	A	283	ASP	CB-CG-OD2	-5.73	113.14	118.30
1	A	81	ASP	CB-CG-OD2	-5.70	113.17	118.30

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Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	A	207	ARG	NE-CZ-NH1	5.55	123.08	120.30
1	A	187	ARG	NE-CZ-NH1	5.54	123.07	120.30
1	A	43	LEU	CB-CG-CD1	-5.38	101.86	111.00
1	A	298	LEU	CB-CG-CD2	-5.35	101.90	111.00
1	A	103	LEU	CB-CG-CD1	-5.32	101.95	111.00
1	A	147	ASP	CB-CG-OD1	5.23	123.00	118.30
1	A	116	LEU	N-CA-C	-5.05	97.35	111.00
1	A	48	ARG	NE-CZ-NH2	-5.05	117.78	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	2430	0	2444	29	0
2	A	335	0	0	16	0
All	All	2765	0	2444	29	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 6.

All (29) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (\AA)	Clash overlap (\AA)
1:A:60:PRO:HD3	2:A:463:HOH:O	1.57	1.05
1:A:207:ARG:HG2	2:A:426:HOH:O	1.78	0.83
1:A:22:LYS:HA	1:A:22:LYS:HE3	1.70	0.72
1:A:108:LYS:HE3	1:A:276:MET:O	1.90	0.71
1:A:9:LYS:HE3	1:A:15:ASN:ND2	2.06	0.69
1:A:43:LEU:HD13	2:A:659:HOH:O	1.94	0.67
1:A:47:HIS:HD2	2:A:348:HOH:O	1.78	0.66
1:A:22:LYS:HE3	1:A:22:LYS:CA	2.23	0.65
1:A:108:LYS:HD3	2:A:557:HOH:O	2.03	0.58
1:A:92:LEU:CD1	1:A:108:LYS:HE2	2.35	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:245:HIS:HB2	1:A:248:GLY:O	2.08	0.53
1:A:304:LYS:HD3	1:A:306:PHE:CZ	2.46	0.51
1:A:178:SER:OG	1:A:182:LYS:HD2	2.11	0.50
1:A:178:SER:HG	1:A:182:LYS:HZ3	1.57	0.49
1:A:9:LYS:HE3	1:A:15:ASN:HD21	1.74	0.48
1:A:142:LYS:NZ	2:A:442:HOH:O	2.46	0.48
1:A:332:GLN:NE2	2:A:459:HOH:O	2.47	0.47
1:A:114:LYS:HB2	2:A:545:HOH:O	2.14	0.47
1:A:108:LYS:CE	2:A:557:HOH:O	2.64	0.46
1:A:20:GLU:OE1	1:A:22:LYS:NZ	2.48	0.46
1:A:146:GLY:HA2	2:A:549:HOH:O	2.15	0.45
1:A:129:LYS:NZ	2:A:461:HOH:O	2.36	0.44
1:A:9:LYS:HE2	1:A:9:LYS:HB3	1.30	0.43
1:A:108:LYS:HA	1:A:108:LYS:HD2	1.60	0.42
1:A:229:LYS:HD2	2:A:555:HOH:O	2.19	0.42
1:A:108:LYS:CD	2:A:557:HOH:O	2.62	0.42
1:A:8:HIS:HA	2:A:658:HOH:O	2.21	0.41
1:A:108:LYS:NZ	2:A:557:HOH:O	2.27	0.41
1:A:277:SER:HA	2:A:411:HOH:O	2.21	0.41

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	318/333 (96%)	311 (98%)	7 (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	255/265 (96%)	252 (99%)	3 (1%)	71 40

All (3) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	9	LYS
1	A	22	LYS
1	A	134	GLN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	47	HIS
1	A	134	GLN
1	A	332	GLN

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	322/333 (96%)	0.42	14 (4%) 35 32	11, 17, 30, 46	0

All (14) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	46	TYR	5.6
1	A	249	VAL	5.0
1	A	60	PRO	3.6
1	A	78	GLY	3.4
1	A	248	GLY	3.0
1	A	197	TYR	2.7
1	A	198	SER	2.6
1	A	49	ALA	2.4
1	A	247	SER	2.4
1	A	245	HIS	2.3
1	A	181	GLU	2.3
1	A	199	THR	2.2
1	A	246	ALA	2.1
1	A	267	ILE	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

There are no ligands in this entry.

6.5 Other polymers

There are no such residues in this entry.