

wwPDB X-ray Structure Validation Summary Report (i)

May 17, 2020 – 02:38 am BST

PDB ID : 3VYH

Title : Crystal structure of aW116R mutant of nitrile hydratase from Pseudonocardia

thermophilla

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Deposited on : 2012-09-25

Resolution : 1.63 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

Mogul: 1.8.5 (274361), CSD as541be (2020)

Xtriage (Phenix) : 1.13 EDS : 2.11

Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)

Refmac: 5.8.0158

 $\begin{array}{cccc} & CCP4 & : & 7.0.044 \; (Gargrove) \\ Ideal \; geometry \; (proteins) & : & Engh \; \& \; Huber \; (2001) \end{array}$

Ideal geometry (DNA, RNA) : Parkinson et al. (1996)

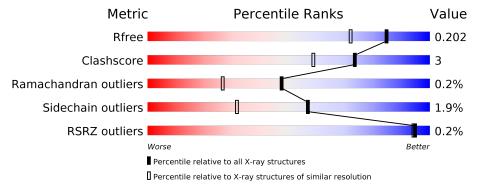
Validation Pipeline (wwPDB-VP) : 2.11

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 1.63 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	$\begin{array}{c} \text{Whole archive} \\ (\#\text{Entries}) \end{array}$	$egin{aligned} ext{Similar resolution} \ (\# ext{Entries, resolution range}(ext{Å})) \end{aligned}$		
R_{free}	130704	3122 (1.66-1.62)		
Clashscore	141614	3268 (1.66-1.62)		
Ramachandran outliers	138981	3215 (1.66-1.62)		
Sidechain outliers	138945	3215 (1.66-1.62)		
RSRZ outliers	127900	3079 (1.66-1.62)		

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	204	92%	7%				
2	В	233	84%	11% • •				



2 Entry composition (i)

There are 4 unique types of molecules in this entry. The entry contains 3818 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 Cobalt-containing nitrile hydratase subunit alpha.

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace		
1	Λ	203	Total	С	N	О	S	0	0	0
1	A	203	1613	1018	273	311	11	0	0	

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
Α	116	ARG	TRP	ENGINEERED MUTATION	UNP Q7SID2

• Molecule 2 is a protein called Cobalt-containing nitrile hydratase subunit beta.

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace		
9	В	226	Total	С	N	О	S	10	0	0
	Б	220	1836	1174	318	338	6	10	0	U

• Molecule 3 is COBALT (II) ION (three-letter code: CO) (formula: Co).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total Co 1 1	0	0

• Molecule 4 is water.

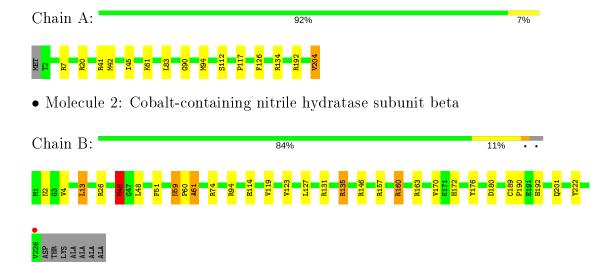
Mo	$1 \mid C$	hain	Residues	${f Atoms}$	ZeroOcc	AltConf
4		A	135	Total O 135 135	0	0
4		В	233	Total O 233 233	0	0



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: Cobalt-containing nitrile hydratase subunit alpha





4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 32 2 1	Depositor
Cell constants	65.68	Depositor
a, b, c, α , β , γ	90.00° 90.00° 120.00°	Depositor
Resolution (Å)	48.44 - 1.63	Depositor
Resolution (A)	48.44 - 1.63	EDS
% Data completeness	99.9 (48.44-1.63)	Depositor
(in resolution range)	99.9 (48.44-1.63)	EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) > 1$	$8.31 \; ({\rm at} \; 1.63 {\rm \AA})$	Xtriage
Refinement program	REFMAC 5.5.0102	Depositor
D.D.	0.174 , 0.203	Depositor
R, R_{free}	0.173 , 0.202	DCC
R_{free} test set	2962 reflections (5.05%)	wwPDB-VP
Wilson B-factor (Å ²)	15.3	Xtriage
Anisotropy	0.002	Xtriage
Bulk solvent $k_{sol}(e/Å^3)$, $B_{sol}(Å^2)$	0.36 , 40.8	EDS
L-test for twinning ²	$< L >=0.47, < L^2>=0.30$	Xtriage
Estimated twinning fraction	0.034 for -h,-k,l	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	3818	wwPDB-VP
Average B, all atoms (Å ²)	16.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: The largest off-origin peak in the Patterson function is 4.39% of the height of the origin peak. No significant pseudotranslation is detected.

²Theoretical 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.



¹Intensities 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: CSD, CSO, CO

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	Bo	nd lengths	Bond angles		
IVIOI	Chain	RMSZ	# Z > 5	RMSZ	# Z > 5	
1	A	1.31	$1/1630 \ (0.1\%)$	1.18	8/2206 (0.4%)	
2	В	1.37	$5/1898 \ (0.3\%)$	1.24	$12/2582 \ (0.5\%)$	
All	All	1.34	$6/3528 \ (0.2\%)$	1.21	$20/4788 \; (0.4\%)$	

The worst 5 of 6 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	$\operatorname{Observed}(\operatorname{\AA})$	$\operatorname{Ideal}(ext{\AA})$
2	В	160	ARG	CG-CD	6.22	1.67	1.51
2	В	176	TYR	CG-CD1	5.91	1.46	1.39
1	A	126	PHE	CE2-CZ	5.38	1.47	1.37
2	В	114	GLU	CD-OE1	5.30	1.31	1.25
2	В	176	TYR	CD2-CE2	5.17	1.47	1.39

The worst 5 of 20 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	\mathbf{Z}	$\mathbf{Observed}(^o)$	$\operatorname{Ideal}(^{o})$
2	В	46	MET	CG-SD-CE	-16.08	74.47	100.20
1	A	134	ARG	NE-CZ-NH2	-10.19	115.21	120.30
2	В	13	LEU	CB-CG-CD1	8.45	125.36	111.00
1	A	7	ARG	NE-CZ-NH1	8.45	124.52	120.30
2	В	135	ARG	NE-CZ-NH2	-7.33	116.64	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)	$\mathbf{H}(\mathbf{added})$	Clashes	Symm-Clashes
1	A	1613	0	1606	7	0
2	В	1836	0	1727	17	0
3	A	1	0	0	0	0
4	A	135	0	0	1	0
4	В	233	0	0	8	1
All	All	3818	0	3333	22	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 22 close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	$egin{array}{c} ext{Interatomic} \ ext{distance } (ext{Å}) \end{array}$	Clash overlap (Å)
1:A:83:LEU:HG	4:A:491:HOH:O	1.66	0.96
2:B:201:GLN:HG2	4:B:340:HOH:O	1.75	0.86
2:B:2:ASN:HD21	2:B:163:ARG:HH12	1.29	0.80
1:A:94:MET:SD	4:B:498:HOH:O	2.49	0.71
2:B:160:ARG:NE	4:B:480:HOH:O	2.30	0.64

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	$egin{array}{c} ext{Interatomic} \ ext{distance } (ext{Å}) \end{array}$	$egin{aligned} ext{Clash} \ ext{overlap } (ext{Å}) \end{aligned}$
4:B:496:HOH:O	4:B:507:HOH:O[1_455]	1.86	0.34

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.

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		* * *				
Mol	Chain	${f Analysed}$	Favoured	Allowed	Outliers	Percentiles
			l	l		l
			T		1	
Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles

Mol	Chain	ain Analysed Favoured Allowed		Outliers	Perce	$_{ m ntiles}$	
1	A	199/204~(98%)	195 (98%)	3 (2%)	1 (0%)	29	11
2	В	$224/233 \ (96\%)$	220 (98%)	4 (2%)	0	100	100
All	All	423/437 (97%)	415 (98%)	7 (2%)	1 (0%)	47	26

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	112	SER

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	hain Analysed Rotameric Outliers		Percentiles		
1	A	177/178 (99%)	176 (99%)	1 (1%)	86 75	
2	В	188/191 (98%)	182 (97%)	6 (3%)	39 12	
All	All	365/369~(99%)	358 (98%)	7 (2%)	57 32	

5 of 7 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	B 59		ASN
2	В	180	ASP
2	В	127	LEU
2	В	13	LEU
2	В	135	ARG

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

Mol	Chain	Res	Type
2	В	2	ASN
2	В	59	ASN

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Mol	Chain	Res	Type
2	В	172	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)

2 non-standard protein/DNA/RNA residues are modelled in this entry.

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).

Mol	Type	Chain	Res	Link	B	ond leng	${ m gths}$	В	ond ang	gles
WIOI	Type	Chain	res	LIIIK	Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
1	CSO	A	113	1,3	3,6,7	1.23	0	0,6,8	0.00	-
1	CSD	A	111	1,3	3,7,8	1.25	0	1,8,10	1.71	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
1	CSO	A	113	1,3	-	1/1/5/7	-
1	CSD	A	111	1,3	-	1/2/6/8	_

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	A	111	CSD	CA-CB-SG-OD1
1	A	113	CSO	N-CA-CB-SG



There are no ring outliers.

No monomer is involved in short contacts.

5.5 Carbohydrates (i)

There are no carbohydrates in this entry.

5.6 Ligand geometry (i)

Of 1 ligands modelled in this entry, 1 is 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, 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 $>$	$\#\mathrm{RSRZ}{>}2$	$\mathbf{OWAB}(\mathrm{\AA}^2)$	Q<0.9
1	A	201/204 (98%)	-0.58	0 100 100	10, 15, 25, 32	0
2	В	$226/233 \ (96\%)$	-0.53	1 (0%) 92 92	9, 13, 23, 30	2 (0%)
All	All	427/437 (97%)	-0.55	1 (0%) 95 94	9, 14, 24, 32	2 (0%)

All (1) RSRZ outliers are listed below:

Mol	Chain	${f Res}$	Type	RSRZ
2	В	226	VAL	2.4

6.2 Non-standard residues in protein, DNA, RNA chains (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	${f B\text{-factors}}({f \AA}^2)$	Q < 0.9
1	CSO	A	113	7/8	0.99	0.06	10,10,13,16	0
1	CSD	A	111	8/9	0.99	0.06	9,9,12,13	0

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. 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	${f B-factors}({f \AA}^2)$	Q < 0.9
3	CO	A	301	1/1	1.00	0.05	10,10,10,10	0

6.5 Other polymers (i)

There are no such residues in this entry.

