



Full wwPDB X-ray Structure Validation Report ⓘ

Sep 29, 2024 – 08:58 AM EDT

PDB ID : 4DVJ
Title : Crystal structure of a putative zinc-dependent alcohol dehydrogenase protein from *Rhizobium etli* CFN 42
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Deposited on : 2012-02-23
Resolution : 1.99 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 types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity	:	4.02b-467
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	1.20.1
EDS	:	3.0
Percentile statistics	:	20231227.v01 (using entries in the PDB archive December 27th 2023)
CCP4	:	9.0.003 (Gargrove)
Density-Fitness	:	1.0.11
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.39

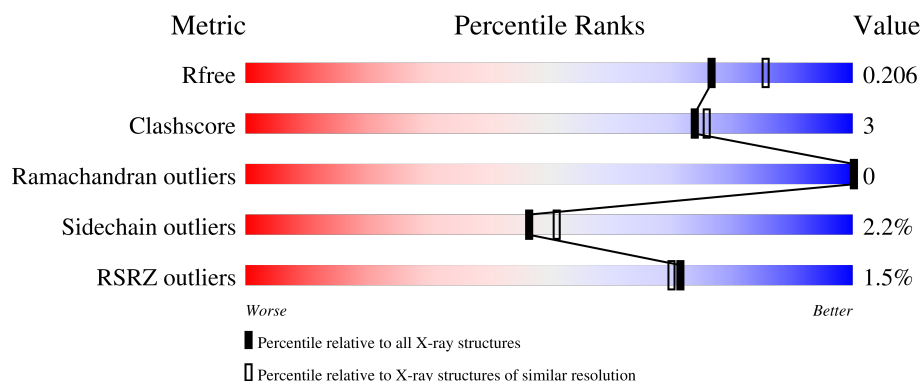
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION



The reported resolution of this entry is 1.99 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}	164625	9409 (2.00-2.00)
Clashscore	180529	10737 (2.00-2.00)
Ramachandran outliers	177936	10628 (2.00-2.00)
Sidechain outliers	177891	10627 (2.00-2.00)
RSRZ outliers	164620	9409 (2.00-2.00)

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 of 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	363	
1	B	363	

2 Entry composition [i](#)

There are 2 unique types of molecules in this entry. The entry contains 5582 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 Putative zinc-dependent alcohol dehydrogenase protein.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	341	Total	C	N	O	S	Se	0	0	0
			2569	1637	452	474	1	5			
1	B	342	Total	C	N	O	S	Se	0	0	0
			2573	1638	452	477	1	5			

There are 44 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	MSE	-	expression tag	UNP Q2JZD9
A	2	HIS	-	expression tag	UNP Q2JZD9
A	3	HIS	-	expression tag	UNP Q2JZD9
A	4	HIS	-	expression tag	UNP Q2JZD9
A	5	HIS	-	expression tag	UNP Q2JZD9
A	6	HIS	-	expression tag	UNP Q2JZD9
A	7	HIS	-	expression tag	UNP Q2JZD9
A	8	SER	-	expression tag	UNP Q2JZD9
A	9	SER	-	expression tag	UNP Q2JZD9
A	10	GLY	-	expression tag	UNP Q2JZD9
A	11	VAL	-	expression tag	UNP Q2JZD9
A	12	ASP	-	expression tag	UNP Q2JZD9
A	13	LEU	-	expression tag	UNP Q2JZD9
A	14	GLY	-	expression tag	UNP Q2JZD9
A	15	THR	-	expression tag	UNP Q2JZD9
A	16	GLU	-	expression tag	UNP Q2JZD9
A	17	ASN	-	expression tag	UNP Q2JZD9
A	18	LEU	-	expression tag	UNP Q2JZD9
A	19	TYR	-	expression tag	UNP Q2JZD9
A	20	PHE	-	expression tag	UNP Q2JZD9
A	21	GLN	-	expression tag	UNP Q2JZD9
A	22	SER	-	expression tag	UNP Q2JZD9
B	1	MSE	-	expression tag	UNP Q2JZD9
B	2	HIS	-	expression tag	UNP Q2JZD9
B	3	HIS	-	expression tag	UNP Q2JZD9

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Chain	Residue	Modelled	Actual	Comment	Reference
B	4	HIS	-	expression tag	UNP Q2JZD9
B	5	HIS	-	expression tag	UNP Q2JZD9
B	6	HIS	-	expression tag	UNP Q2JZD9
B	7	HIS	-	expression tag	UNP Q2JZD9
B	8	SER	-	expression tag	UNP Q2JZD9
B	9	SER	-	expression tag	UNP Q2JZD9
B	10	GLY	-	expression tag	UNP Q2JZD9
B	11	VAL	-	expression tag	UNP Q2JZD9
B	12	ASP	-	expression tag	UNP Q2JZD9
B	13	LEU	-	expression tag	UNP Q2JZD9
B	14	GLY	-	expression tag	UNP Q2JZD9
B	15	THR	-	expression tag	UNP Q2JZD9
B	16	GLU	-	expression tag	UNP Q2JZD9
B	17	ASN	-	expression tag	UNP Q2JZD9
B	18	LEU	-	expression tag	UNP Q2JZD9
B	19	TYR	-	expression tag	UNP Q2JZD9
B	20	PHE	-	expression tag	UNP Q2JZD9
B	21	GLN	-	expression tag	UNP Q2JZD9
B	22	SER	-	expression tag	UNP Q2JZD9

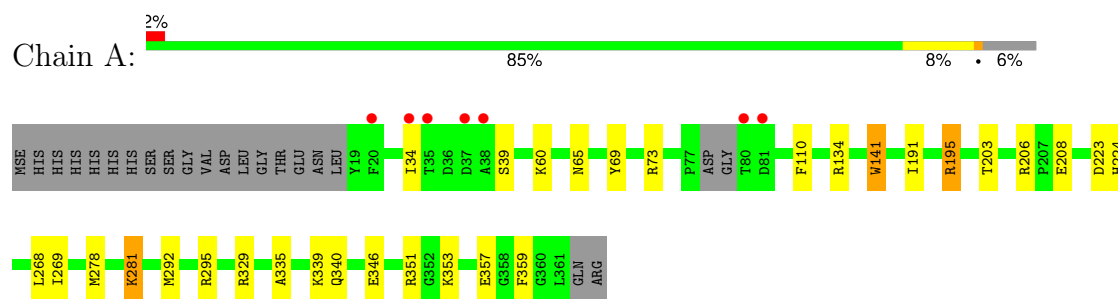
- Molecule 2 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	203	Total O 203 203	0	0
2	B	237	Total O 237 237	0	0

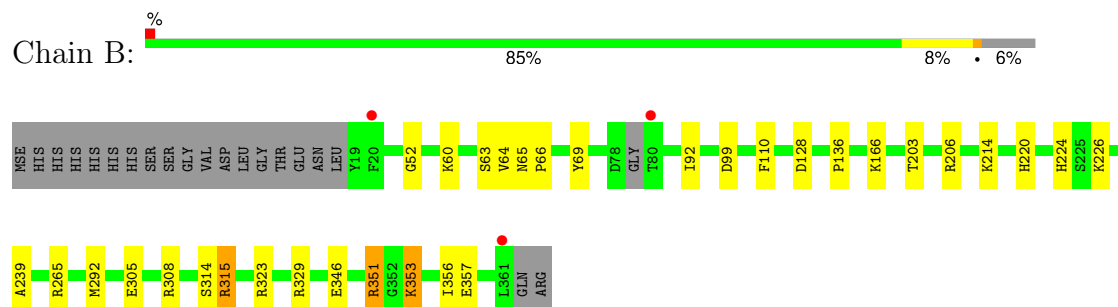
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide 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: Putative zinc-dependent alcohol dehydrogenase protein



- Molecule 1: Putative zinc-dependent alcohol dehydrogenase protein



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	185.31Å 42.24Å 111.27Å 90.00° 122.39° 90.00°	Depositor
Resolution (Å)	48.66 – 1.99 48.66 – 1.99	Depositor EDS
% Data completeness (in resolution range)	99.5 (48.66-1.99) 99.5 (48.66-1.99)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	5.01 (at 1.98Å)	Xtriage
Refinement program	REFMAC 5.5.0109	Depositor
R, R_{free}	0.162 , 0.206 0.164 , 0.206	Depositor DCC
R_{free} test set	2573 reflections (5.08%)	wwPDB-VP
Wilson B-factor (Å ²)	21.6	Xtriage
Anisotropy	0.037	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 32.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	5582	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 5.38% 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

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	1.21	1/2622 (0.0%)	0.96	2/3565 (0.1%)
1	B	1.27	3/2626 (0.1%)	0.98	7/3572 (0.2%)
All	All	1.24	4/5248 (0.1%)	0.97	9/7137 (0.1%)

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	64	VAL	CB-CG1	5.84	1.65	1.52
1	B	239	ALA	CA-CB	5.74	1.64	1.52
1	B	351	ARG	CB-CG	-5.29	1.38	1.52
1	A	141	TRP	CE3-CZ3	5.20	1.47	1.38

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	128	ASP	CB-CG-OD2	-5.84	113.04	118.30
1	B	206	ARG	NE-CZ-NH2	-5.67	117.47	120.30
1	B	353	LYS	CD-CE-NZ	5.47	124.28	111.70
1	A	223	ASP	CB-CG-OD1	5.32	123.09	118.30
1	B	265	ARG	NE-CZ-NH2	-5.30	117.65	120.30
1	B	265	ARG	NE-CZ-NH1	5.21	122.91	120.30
1	A	295	ARG	NE-CZ-NH2	-5.11	117.74	120.30
1	B	315	ARG	NE-CZ-NH1	5.03	122.82	120.30
1	B	226	LYS	CD-CE-NZ	-5.03	100.14	111.70

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	2569	0	2585	16	0
1	B	2573	0	2578	15	0
2	A	203	0	0	3	0
2	B	237	0	0	6	0
All	All	5582	0	5163	31	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 3.

All (31) 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:203:THR:OG1	1:A:224:HIS:HD2	1.70	0.75
1:A:278:MSE:HE1	1:A:281:LYS:HE2	1.74	0.69
1:B:214:LYS:HE3	2:B:579:HOH:O	1.94	0.67
1:B:220:HIS:HE1	2:B:634:HOH:O	1.78	0.66
1:A:65:ASN:OD1	1:A:353:LYS:NZ	2.30	0.63
1:B:203:THR:OG1	1:B:224:HIS:HD2	1.80	0.63
1:B:60:LYS:HE2	1:B:92:ILE:HD12	1.85	0.58
1:B:329:ARG:HD3	1:B:357:GLU:OE2	2.04	0.58
1:B:305:GLU:OE2	1:B:308:ARG:NH1	2.41	0.54
1:A:134:ARG:HG2	2:A:541:HOH:O	2.09	0.51
1:A:206:ARG:HE	1:A:208:GLU:CD	2.15	0.50
1:A:359:PHE:HA	2:A:520:HOH:O	2.12	0.50
1:B:292:MSE:HG3	2:B:546:HOH:O	2.12	0.49
1:A:60:LYS:HG3	1:A:141:TRP:CD2	2.48	0.48
1:A:329:ARG:HD3	1:A:357:GLU:OE2	2.14	0.47
1:B:65:ASN:OD1	1:B:353:LYS:NZ	2.47	0.47
1:B:356:ILE:HD12	1:B:356:ILE:N	2.30	0.47
1:B:136:PRO:HG3	1:B:314:SER:HA	1.96	0.47
1:A:335:ALA:O	1:A:339:LYS:HG3	2.15	0.47
1:B:323:ARG:HB2	2:B:460:HOH:O	2.14	0.46
1:B:69:TYR:CG	1:B:346:GLU:HG2	2.51	0.46
1:A:195:ARG:NE	1:A:195:ARG:HA	2.30	0.46
1:A:69:TYR:O	1:A:73:ARG:HG3	2.15	0.45
1:B:315:ARG:NH2	2:B:597:HOH:O	2.50	0.43
1:A:191:ILE:O	1:A:195:ARG:HB2	2.20	0.42
1:A:203:THR:OG1	1:A:224:HIS:CD2	2.61	0.41
1:A:34:ILE:HD12	1:A:39:SER:OG	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:315:ARG:NH2	2:B:633:HOH:O	2.53	0.41
1:B:52:GLY:HA2	1:B:99:ASP:OD2	2.21	0.40
1:A:69:TYR:CG	1:A:346:GLU:HG2	2.56	0.40
1:A:292:MSE:HG3	2:A:418:HOH:O	2.22	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	337/363 (93%)	329 (98%)	8 (2%)	0	100	100
1	B	338/363 (93%)	329 (97%)	9 (3%)	0	100	100
All	All	675/726 (93%)	658 (98%)	17 (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	268/285 (94%)	261 (97%)	7 (3%)	41	44
1	B	268/285 (94%)	263 (98%)	5 (2%)	52	57
All	All	536/570 (94%)	524 (98%)	12 (2%)	47	51

All (12) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	110	PHE
1	A	195	ARG
1	A	268	LEU
1	A	269	ILE
1	A	281	LYS
1	A	340	GLN
1	A	351	ARG
1	B	63	SER
1	B	66	PRO
1	B	110	PHE
1	B	166	LYS
1	B	351	ARG

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

Mol	Chain	Res	Type
1	A	224	HIS
1	B	220	HIS
1	B	224	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 oligosaccharides 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	336/363 (92%)	-0.40	7 (2%) 63 62	12, 21, 38, 58	0
1	B	337/363 (92%)	-0.57	3 (0%) 81 80	10, 19, 33, 44	0
All	All	673/726 (92%)	-0.49	10 (1%) 71 70	10, 20, 35, 58	0

All (10) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	37	ASP	4.4
1	A	38	ALA	3.9
1	B	80	THR	3.3
1	A	35	THR	3.2
1	A	20	PHE	3.2
1	B	20	PHE	3.1
1	A	80	THR	3.1
1	A	34	ILE	2.8
1	A	81	ASP	2.1
1	B	361	LEU	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 monosaccharides 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.