



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 03:45 PM UTC

PDB ID : 5UX1 / pdb_00005ux1
Title : Protein 43 with aldehyde deformylating oxygenase activity from *Synechococcus*
Authors : Wilson, D.K.; Mak, W.S.; Siegel, J.B.
Deposited on : 2017-02-21
Resolution : 1.50 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	:	4-5-2 with Phenix2.0
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

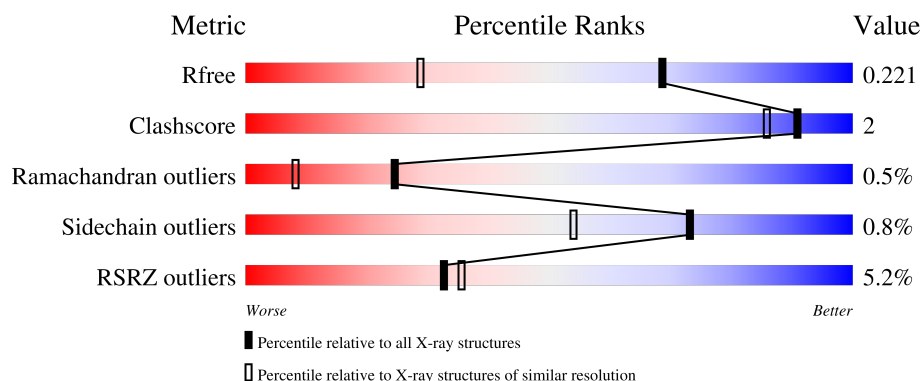
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.50 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}	180053	4037 (1.50-1.50)
Clashscore	190562	4235 (1.50-1.50)
Ramachandran outliers	187476	4153 (1.50-1.50)
Sidechain outliers	187428	4150 (1.50-1.50)
RSRZ outliers	180081	4039 (1.50-1.50)

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	237	3% 80% 5% 15%
1	B	237	2% 82% • 15%
1	C	237	4% 84% • 15%
1	D	237	8% 75% 8% 17%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 6905 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 tRNA-(MS(2)IO(6)A)-hydroxylase-like.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	201	Total	C	N	O	S	0	0	0
			1599	1012	288	289	10			
1	B	201	Total	C	N	O	S	0	0	0
			1599	1012	288	289	10			
1	C	202	Total	C	N	O	S	0	0	0
			1605	1015	289	291	10			
1	D	197	Total	C	N	O	S	0	0	0
			1564	991	280	283	10			

There are 32 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	230	GLY	-	expression tag	UNP Q3AMA5
A	231	SER	-	expression tag	UNP Q3AMA5
A	232	HIS	-	expression tag	UNP Q3AMA5
A	233	HIS	-	expression tag	UNP Q3AMA5
A	234	HIS	-	expression tag	UNP Q3AMA5
A	235	HIS	-	expression tag	UNP Q3AMA5
A	236	HIS	-	expression tag	UNP Q3AMA5
A	237	HIS	-	expression tag	UNP Q3AMA5
B	230	GLY	-	expression tag	UNP Q3AMA5
B	231	SER	-	expression tag	UNP Q3AMA5
B	232	HIS	-	expression tag	UNP Q3AMA5
B	233	HIS	-	expression tag	UNP Q3AMA5
B	234	HIS	-	expression tag	UNP Q3AMA5
B	235	HIS	-	expression tag	UNP Q3AMA5
B	236	HIS	-	expression tag	UNP Q3AMA5
B	237	HIS	-	expression tag	UNP Q3AMA5
C	230	GLY	-	expression tag	UNP Q3AMA5
C	231	SER	-	expression tag	UNP Q3AMA5
C	232	HIS	-	expression tag	UNP Q3AMA5
C	233	HIS	-	expression tag	UNP Q3AMA5
C	234	HIS	-	expression tag	UNP Q3AMA5

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Chain	Residue	Modelled	Actual	Comment	Reference
C	235	HIS	-	expression tag	UNP Q3AMA5
C	236	HIS	-	expression tag	UNP Q3AMA5
C	237	HIS	-	expression tag	UNP Q3AMA5
D	230	GLY	-	expression tag	UNP Q3AMA5
D	231	SER	-	expression tag	UNP Q3AMA5
D	232	HIS	-	expression tag	UNP Q3AMA5
D	233	HIS	-	expression tag	UNP Q3AMA5
D	234	HIS	-	expression tag	UNP Q3AMA5
D	235	HIS	-	expression tag	UNP Q3AMA5
D	236	HIS	-	expression tag	UNP Q3AMA5
D	237	HIS	-	expression tag	UNP Q3AMA5

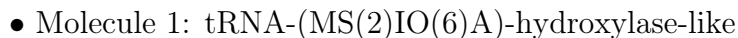
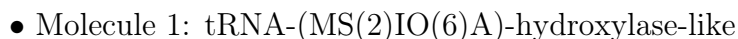
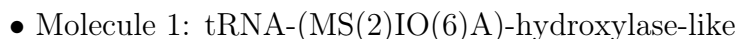
- Molecule 2 is FE (III) ION (CCD ID: FE) (formula: Fe).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	2	Total Fe 2 2	0	2
2	B	2	Total Fe 2 2	0	2
2	C	2	Total Fe 2 2	0	2
2	D	2	Total Fe 2 2	0	2

- Molecule 3 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	131	Total O 131 131	0	0
3	B	149	Total O 149 149	0	0
3	C	164	Total O 164 164	0	0
3	D	86	Total O 86 86	0	0

- Molecule 1: tRNA-(MS(2)IO(6)A)-hydroxylase-like



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	72.34Å 78.27Å 83.73Å 90.00° 103.20° 90.00°	Depositor
Resolution (Å)	81.50 – 1.50 81.50 – 1.50	Depositor EDS
% Data completeness (in resolution range)	99.3 (81.50-1.50) 99.4 (81.50-1.50)	Depositor EDS
R_{merge}	0.04	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.25 (at 1.50Å)	Xtriage
Refinement program	REFMAC 5.8.0049	Depositor
R, R_{free}	0.192 , 0.215 (Not available) , 0.221	Depositor DCC
R_{free} test set	3619 reflections (2.50%)	wwPDB-VP
Wilson B-factor (Å ²)	18.6	Xtriage
Anisotropy	0.047	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 26.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	6905	wwPDB-VP
Average B, all atoms (Å ²)	23.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.65% 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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: FE

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	0.74	0/1636	0.85	0/2218
1	B	0.70	0/1636	0.84	0/2218
1	C	0.78	0/1642	0.89	0/2226
1	D	0.65	0/1601	0.82	1/2171 (0.0%)
All	All	0.72	0/6515	0.85	1/8833 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	120	LYS	N-CA-C	5.79	119.53	112.23

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	1599	0	1587	6	0
1	B	1599	0	1587	4	0
1	C	1605	0	1592	2	0
1	D	1564	0	1548	9	0
2	A	2	0	0	0	0
2	B	2	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	C	2	0	0	0	0
2	D	2	0	0	0	0
3	A	131	0	0	0	0
3	B	149	0	0	1	0
3	C	164	0	0	0	0
3	D	86	0	0	0	0
All	All	6905	0	6314	21	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 2.

All (21) 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:113:ASP:OD1	1:A:116:ARG:NH2	2.20	0.75
1:D:185:GLU:H	1:D:185:GLU:CD	2.16	0.54
1:A:84:GLU:HA	1:A:84:GLU:OE1	2.09	0.53
1:B:80:LEU:C	1:B:80:LEU:HD23	2.35	0.51
1:A:46:LEU:HD22	1:A:91:VAL:HG13	1.93	0.51
1:A:168:ALA:HA	1:A:171:PHE:CE2	2.48	0.49
1:B:46:LEU:HD22	1:B:91:VAL:HG13	1.94	0.48
1:D:66:ARG:HD2	1:D:67:TYR:CZ	2.48	0.48
1:D:98:ARG:NH2	1:D:153:ASP:OD2	2.41	0.48
1:D:46:LEU:HD22	1:D:91:VAL:HG13	1.96	0.47
1:D:144:MET:HE1	1:D:167:GLU:HG3	1.97	0.46
1:D:43:MET:HE2	1:D:100:ARG:HG3	1.99	0.45
1:D:168:ALA:HA	1:D:171:PHE:CE2	2.53	0.44
1:B:168:ALA:HA	1:B:171:PHE:CE2	2.53	0.44
1:D:84:GLU:HA	1:D:84:GLU:OE1	2.18	0.43
1:C:144:MET:HE1	1:C:167:GLU:HG3	2.01	0.43
1:D:83:GLU:CD	1:D:170:HIS:HE2	2.27	0.43
1:C:84:GLU:HA	1:C:84:GLU:OE1	2.19	0.42
1:A:80:LEU:C	1:A:80:LEU:HD23	2.45	0.41
1:A:92:LEU:HD13	1:A:102:LEU:HG	2.03	0.41
1:B:66:ARG:HD2	3:B:456:HOH:O	2.20	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	199/237 (84%)	197 (99%)	1 (0%)	1 (0%)	24	8
1	B	199/237 (84%)	197 (99%)	1 (0%)	1 (0%)	24	8
1	C	200/237 (84%)	198 (99%)	1 (0%)	1 (0%)	24	8
1	D	195/237 (82%)	192 (98%)	2 (1%)	1 (0%)	24	8
All	All	793/948 (84%)	784 (99%)	5 (1%)	4 (0%)	24	8

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	218	GLY
1	A	218	GLY
1	B	218	GLY
1	C	218	GLY

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	165/198 (83%)	163 (99%)	2 (1%)	63	38
1	B	165/198 (83%)	164 (99%)	1 (1%)	78	62
1	C	166/198 (84%)	166 (100%)	0	100	100
1	D	161/198 (81%)	159 (99%)	2 (1%)	63	38
All	All	657/792 (83%)	652 (99%)	5 (1%)	73	54

All (5) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	46	LEU
1	A	120	LYS
1	B	46	LEU
1	D	124	GLN
1	D	179	GLU

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	117	GLN
1	B	117	GLN
1	D	124	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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 8 ligands modelled in this entry, 8 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

6.1 Protein, DNA and RNA chains

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	201/237 (84%)	0.33	8 (3%) 42 46	13, 20, 34, 48	0
1	B	201/237 (84%)	0.22	4 (1%) 65 68	14, 20, 32, 42	0
1	C	202/237 (85%)	0.07	10 (4%) 34 37	12, 17, 29, 41	0
1	D	197/237 (83%)	0.80	20 (10%) 12 12	13, 27, 41, 50	0
All	All	801/948 (84%)	0.35	42 (5%) 33 35	12, 21, 36, 50	0

All (42) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	221	VAL	5.1
1	A	23	ILE	4.6
1	C	224	ILE	4.1
1	D	207	THR	3.8
1	C	222	THR	3.1
1	A	223	GLN	3.0
1	C	225	SER	2.7
1	C	221	VAL	2.7
1	D	193	VAL	2.7
1	D	186	LEU	2.6
1	D	32	TRP	2.6
1	D	99	GLY	2.6
1	C	24	ARG	2.6
1	B	23	ILE	2.5
1	D	27	ALA	2.5
1	C	69	CYS	2.5
1	A	162	ASP	2.5
1	A	163	LEU	2.4
1	D	114	LEU	2.4
1	D	97	ALA	2.4
1	A	165	ALA	2.3

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Mol	Chain	Res	Type	RSRZ
1	C	223	GLN	2.3
1	D	206	LEU	2.3
1	D	179	GLU	2.3
1	B	223	GLN	2.2
1	D	120	LYS	2.2
1	A	164	LEU	2.2
1	C	186	LEU	2.2
1	D	118	ILE	2.2
1	D	161	SER	2.2
1	B	162	ASP	2.2
1	A	167	GLU	2.2
1	D	209	PRO	2.1
1	D	40	ALA	2.1
1	C	120	LYS	2.0
1	D	115	ALA	2.0
1	D	196	LEU	2.0
1	A	185	GLU	2.0
1	C	162	ASP	2.0
1	D	98	ARG	2.0
1	B	32	TRP	2.0
1	D	105	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 oligosaccharides 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, 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	B-factors($\text{\AA}^2$)	Q<0.9
2	FE	A	302[L]	1/1	0.97	0.04	24,24,24,24	1
2	FE	B	302[L]	1/1	0.97	0.04	22,22,22,22	1

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	FE	A	301[L]	1/1	0.98	0.03	23,23,23,23	1
2	FE	C	301[L]	1/1	0.98	0.04	19,19,19,19	1
2	FE	C	302[L]	1/1	0.98	0.03	22,22,22,22	1
2	FE	D	301[L]	1/1	0.98	0.04	25,25,25,25	1
2	FE	D	302[L]	1/1	0.98	0.05	27,27,27,27	1
2	FE	B	301[L]	1/1	0.99	0.04	22,22,22,22	1

6.5 Other polymers [i](#)

There are no such residues in this entry.