



wwPDB X-ray Structure Validation Summary Report ⓘ

Feb 14, 2017 – 10:44 pm GMT

PDB ID : 2Z1B
Title : Crystal Structure of 5-aminolevulinic acid dehydratase (ALAD) from Mus musculus
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Deposited on : 2007-05-08
Resolution : 3.30 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

<http://wwpdb.org/validation/2016/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
Xtriage (Phenix) : 1.9-1692
EDS : trunk28620
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : recalc28949

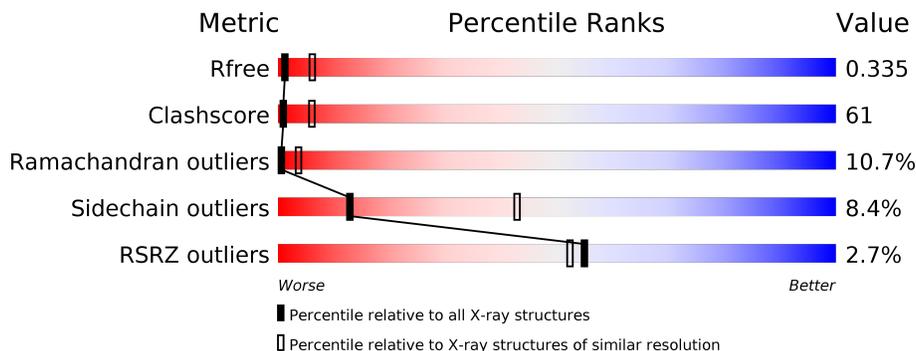
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.30 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}	100719	1034 (3.36-3.24)
Clashscore	112137	1100 (3.36-3.24)
Ramachandran outliers	110173	1081 (3.36-3.24)
Sidechain outliers	110143	1080 (3.36-3.24)
RSRZ outliers	101464	1039 (3.36-3.24)

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. 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	330	 5% 29% 49% 10% 12%
1	B	330	 5% 27% 56% 14% 10%
1	C	330	 5% 20% 54% 13% 13%
1	D	330	 3% 22% 55% 12% 10%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 9223 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 Delta-aminolevulinic acid dehydratase.

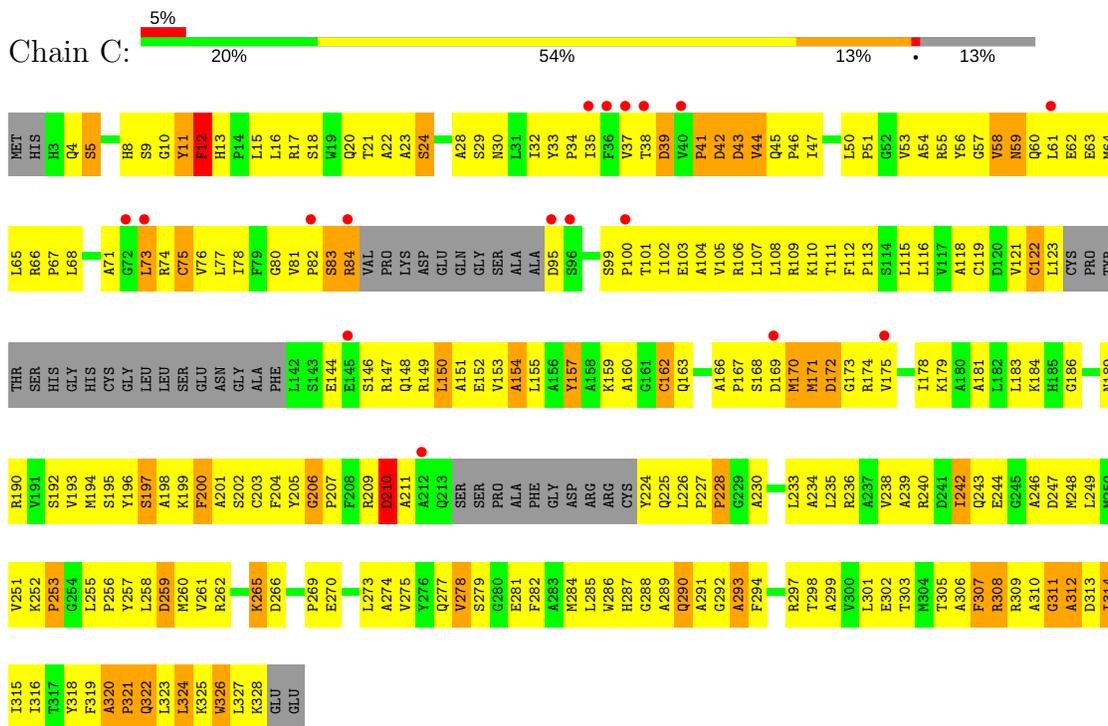
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	290	2230	1428	387	400	15	0	0	0
1	B	320	2445	1557	426	445	17	0	0	0
1	C	288	2217	1420	385	398	14	0	0	0
1	D	296	2274	1454	397	408	15	0	0	0

- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	12	Total	O	0	0
			12	12		
2	B	20	Total	O	0	0
			20	20		
2	C	15	Total	O	0	0
			15	15		
2	D	10	Total	O	0	0
			10	10		

L327
K328
GLU
GLU

● Molecule 1: Delta-aminolevulinic acid dehydratase



4 Data and refinement statistics

Property	Value	Source
Space group	P 62 2 2	Depositor
Cell constants a, b, c, α , β , γ	240.92Å 240.92Å 103.94Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	43.40 – 3.30 47.72 – 3.29	Depositor EDS
% Data completeness (in resolution range)	88.7 (43.40-3.30) 95.4 (47.72-3.29)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.11	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.79 (at 3.33Å)	Xtrriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.278 , 0.353 0.266 , 0.335	Depositor DCC
R_{free} test set	2400 reflections (9.94%)	DCC
Wilson B-factor (Å ²)	98.3	Xtrriage
Anisotropy	0.108	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 90.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.90	EDS
Total number of atoms	9223	wwPDB-VP
Average B, all atoms (Å ²)	70.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.26% 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](#)

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	0.33	0/2278	0.57	0/3091
1	B	0.34	0/2500	0.61	0/3394
1	C	0.31	0/2264	0.57	0/3071
1	D	0.32	0/2323	0.56	0/3151
All	All	0.32	0/9365	0.58	0/12707

There are no bond length outliers.

There are no bond angle outliers.

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	2230	0	2252	229	0
1	B	2445	0	2452	303	0
1	C	2217	0	2240	352	0
1	D	2274	0	2297	324	0
2	A	12	0	0	3	0
2	B	20	0	0	5	0
2	C	15	0	0	4	0
2	D	10	0	0	5	0
All	All	9223	0	9241	1127	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 61.

The worst 5 of 1127 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:227:PRO:HG2	1:A:230:ALA:HB2	1.35	1.09
1:A:5:SER:HB2	1:B:240:ARG:HG3	1.39	1.03
1:A:242:ILE:HD13	1:A:271:LEU:HD11	1.40	1.03
1:C:227:PRO:HG2	1:C:230:ALA:HB2	1.40	1.03
1:C:255:LEU:HD11	1:D:256:PRO:HA	1.40	1.02

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	282/330 (86%)	196 (70%)	68 (24%)	18 (6%)	1	12
1	B	316/330 (96%)	203 (64%)	80 (25%)	33 (10%)	0	4
1	C	280/330 (85%)	179 (64%)	60 (21%)	41 (15%)	0	1
1	D	290/330 (88%)	178 (61%)	79 (27%)	33 (11%)	0	3
All	All	1168/1320 (88%)	756 (65%)	287 (25%)	125 (11%)	0	4

5 of 125 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	58	VAL
1	A	226	LEU
1	A	278	VAL
1	A	290	GLN
1	B	12	PHE

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	235/267 (88%)	214 (91%)	21 (9%)	11	39
1	B	258/267 (97%)	235 (91%)	23 (9%)	11	39
1	C	233/267 (87%)	216 (93%)	17 (7%)	16	49
1	D	239/267 (90%)	219 (92%)	20 (8%)	13	43
All	All	965/1068 (90%)	884 (92%)	81 (8%)	13	43

5 of 81 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	248	MET
1	C	12	PHE
1	D	255	LEU
1	B	255	LEU
1	B	323	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 15 such sidechains are listed below:

Mol	Chain	Res	Type
1	B	59	ASN
1	B	60	GLN
1	D	60	GLN
1	B	45	GLN
1	D	59	ASN

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	290/330 (87%)	-0.09	2 (0%) 87 87	25, 54, 106, 135	0
1	B	320/330 (96%)	-0.08	4 (1%) 77 75	24, 56, 101, 127	0
1	C	288/330 (87%)	0.19	17 (5%) 23 22	36, 80, 126, 148	0
1	D	296/330 (89%)	0.18	9 (3%) 51 49	34, 83, 130, 143	0
All	All	1194/1320 (90%)	0.05	32 (2%) 55 52	24, 66, 120, 148	0

The worst 5 of 32 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	96	SER	6.2
1	C	95	ASP	5.1
1	C	38	THR	4.5
1	B	131	HIS	3.9
1	C	169	ASP	3.8

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.