



Full wwPDB X-ray Structure Validation Report ⓘ

Oct 12, 2020 – 10:31 AM EDT

PDB ID : 4EPK
Title : Evidence for a Dual Role of an Active Site Histidine in alpha-Amino-beta-Carboxymuconate-epsilon-Semialdehyde Decarboxylase
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Deposited on : 2012-04-17
Resolution : 2.60 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
Xtriage (Phenix)	:	1.13
EDS	:	2.14.6
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)	:	2.14.6

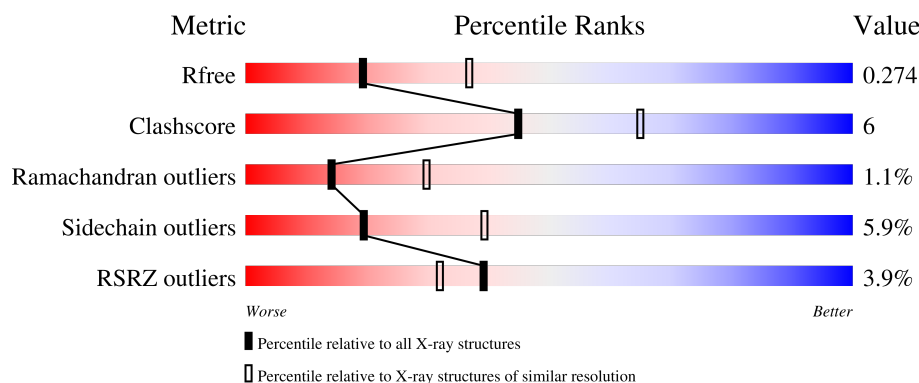
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.60 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	3163 (2.60-2.60)
Clashscore	141614	3518 (2.60-2.60)
Ramachandran outliers	138981	3455 (2.60-2.60)
Sidechain outliers	138945	3455 (2.60-2.60)
RSRZ outliers	127900	3104 (2.60-2.60)

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	334	 83% 15% ..
1	B	334	 78% 19% ..

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 5188 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-amino-3-carboxymuconate 6-semialdehyde decarboxylase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	331	Total	C	N	O	S	1	0	0
			2581	1646	448	469	18			
1	B	330	Total	C	N	O	S	3	0	0
			2572	1640	446	468	18			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	228	GLY	HIS	engineered mutation	UNP Q83V25
B	228	GLY	HIS	engineered mutation	UNP Q83V25

- Molecule 2 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	1	Total	Zn	0	0
			1	1		
2	A	1	Total	Zn	0	0
			1	1		

- Molecule 3 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

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

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	24	Total	O	0	0
			24	24		
4	B	8	Total	O	0	0
			8	8		

- Molecule 1: 2-amino-3-carboxymuconate 6-semialdehyde decarboxylase



4 Data and refinement statistics

Property	Value	Source
Space group	P 42 21 2	Depositor
Cell constants a, b, c, α , β , γ	91.50Å 91.50Å 170.12Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	32.35 – 2.60 32.35 – 2.60	Depositor EDS
% Data completeness (in resolution range)	99.6 (32.35-2.60) 99.6 (32.35-2.60)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.76 (at 2.61Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.7.3_928)	Depositor
R, R_{free}	0.208 , 0.281 0.203 , 0.274	Depositor DCC
R_{free} test set	1173 reflections (5.13%)	wwPDB-VP
Wilson B-factor (Å ²)	65.4	Xtriage
Anisotropy	0.136	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 41.3	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	0.95	EDS
Total number of atoms	5188	wwPDB-VP
Average B, all atoms (Å ²)	68.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.93% 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

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

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	0.47	0/2647	0.60	0/3586
1	B	0.43	0/2638	0.59	1/3574 (0.0%)
All	All	0.45	0/5285	0.59	1/7160 (0.0%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	B	0	1

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	B	155	LEU	CA-CB-CG	5.52	128.00	115.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	154	ASP	Peptide

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	2581	0	2528	32	0
1	B	2572	0	2516	35	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	1	0	0	0	0
4	A	24	0	0	4	0
4	B	8	0	0	1	0
All	All	5188	0	5044	66	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 (66) 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:B:150:LEU:HB2	1:B:153:LYS:HB2	1.42	1.01
1:A:15:ARG:NH2	4:A:523:HOH:O	2.05	0.87
1:B:149:HIS:CD2	1:B:179:TRP:HB2	2.23	0.72
1:A:271:VAL:O	4:A:502:HOH:O	2.07	0.70
1:A:14:PRO:HG2	1:A:16:ILE:HD11	1.75	0.68
1:A:248:ASP:OD1	4:A:521:HOH:O	2.13	0.67
1:A:112:PRO:O	1:A:116:LYS:NZ	2.28	0.65
1:B:37:LYS:O	1:B:39:ASP:N	2.28	0.64
1:B:127:ASP:N	1:B:127:ASP:OD1	2.26	0.62
1:B:150:LEU:HD13	1:B:153:LYS:HD2	1.82	0.60
1:A:256:ARG:NH2	4:A:513:HOH:O	2.34	0.60
1:B:40:THR:OG1	1:B:41:GLY:N	2.27	0.59
1:B:124:GLN:HA	1:B:150:LEU:HD21	1.83	0.59
1:A:18:GLU:HB2	1:A:31:TRP:CZ2	2.41	0.55
1:B:5:ARG:HG3	1:B:322:ILE:HG23	1.89	0.54
1:B:122:PRO:HD2	1:B:129:ALA:HA	1.89	0.54
1:B:48:ASN:N	1:B:48:ASN:OD1	2.41	0.54
1:B:53:VAL:HG12	1:B:58:TRP:HZ3	1.71	0.54
1:A:6:ILE:HG12	1:A:74:VAL:HB	1.90	0.53
1:B:51:ARG:HG2	1:B:52:PRO:HD2	1.89	0.53
1:B:218:ILE:HG21	1:B:222:LEU:HD22	1.91	0.52
1:B:55:GLN:HA	1:B:58:TRP:CH2	2.45	0.52
1:A:18:GLU:HB2	1:A:31:TRP:CE2	2.45	0.51
1:B:156:ASP:OD2	4:B:508:HOH:O	2.19	0.51
1:A:5:ARG:HD3	1:A:71:GLY:O	2.11	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:96:TRP:O	1:B:100:MET:HG2	2.12	0.50
1:B:318:LYS:O	1:B:322:ILE:HG12	2.12	0.49
1:B:137:VAL:HG21	1:B:172:ILE:HG12	1.94	0.49
1:A:327:SER:O	1:A:331:ASN:HA	2.12	0.49
1:A:256:ARG:HB3	1:A:260:GLU:OE1	2.12	0.49
1:A:189:LYS:HB3	1:A:190:TRP:CD2	2.48	0.49
1:A:212:SER:HA	1:B:190:TRP:CZ2	2.47	0.49
1:A:78:CYS:HB3	1:A:118:LEU:HB2	1.97	0.47
1:A:286:GLU:HG3	1:A:320:LYS:HD2	1.96	0.47
1:A:148:ASN:OD1	1:A:178:PRO:HA	2.14	0.46
1:B:214:ALA:O	1:B:218:ILE:HG12	2.16	0.45
1:B:153:LYS:HG2	1:B:154:ASP:H	1.82	0.45
1:B:153:LYS:HG2	1:B:154:ASP:N	2.32	0.45
1:B:156:ASP:OD1	1:B:160:LEU:HD22	2.17	0.45
1:B:203:GLN:HG3	1:B:233:PHE:HA	1.98	0.44
1:B:147:GLY:HA2	1:B:177:HIS:O	2.17	0.44
1:A:172:ILE:HA	1:A:173:PRO:HD3	1.85	0.43
1:B:5:ARG:HD2	1:B:71:GLY:O	2.18	0.43
1:A:220:LYS:HA	1:A:264:ARG:NH2	2.34	0.43
1:A:188:LYS:HE2	1:A:188:LYS:HB3	1.83	0.43
1:B:125:ASP:OD2	1:B:128:LEU:HG	2.19	0.43
1:B:155:LEU:HD13	1:B:156:ASP:N	2.33	0.43
1:A:248:ASP:OD1	1:A:248:ASP:N	2.49	0.42
1:A:37:LYS:NZ	1:A:37:LYS:HB3	2.34	0.42
1:B:42:SER:HB2	1:B:51:ARG:O	2.19	0.42
1:A:175:LEU:HD12	1:A:225:CYS:O	2.19	0.42
1:A:187:MET:HE2	1:A:193:PRO:HA	2.00	0.42
1:B:8:MET:HA	1:B:76:VAL:HB	2.02	0.42
1:A:215:PHE:O	1:A:264:ARG:NH1	2.48	0.42
1:B:11:HIS:O	1:B:63:ARG:NH2	2.51	0.42
1:A:31:TRP:CE2	1:A:44:MET:HB2	2.54	0.42
1:B:218:ILE:HA	1:B:219:PRO:HD3	1.95	0.41
1:A:142:LEU:HA	1:A:142:LEU:HD23	1.84	0.41
1:B:105:LEU:HD23	1:B:105:LEU:HA	1.95	0.41
1:A:257:PRO:HA	1:A:258:PRO:HD3	1.85	0.41
1:B:88:TRP:HB2	1:B:93:ALA:HB2	2.02	0.41
1:A:128:LEU:HD23	1:A:128:LEU:HA	1.85	0.41
1:A:15:ARG:HE	1:A:15:ARG:HB3	1.30	0.41
1:A:188:LYS:HG2	1:A:189:LYS:N	2.35	0.41
1:A:297:PHE:HA	1:A:298:PRO:HD3	1.89	0.41
1:B:218:ILE:O	1:B:264:ARG:NH2	2.50	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	329/334 (98%)	301 (92%)	26 (8%)	2 (1%)	25	47
1	B	328/334 (98%)	293 (89%)	30 (9%)	5 (2%)	10	21
All	All	657/668 (98%)	594 (90%)	56 (8%)	7 (1%)	14	30

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	38	GLY
1	B	270	ALA
1	A	214	ALA
1	B	220	LYS
1	B	59	ASP
1	A	255	PRO
1	B	298	PRO

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	270/273 (99%)	257 (95%)	13 (5%)	25	49
1	B	269/273 (98%)	250 (93%)	19 (7%)	14	29
All	All	539/546 (99%)	507 (94%)	32 (6%)	19	39

All (32) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	15	ARG
1	A	16	ILE
1	A	19	GLN
1	A	37	LYS
1	A	51	ARG
1	A	150	LEU
1	A	179	TRP
1	A	188	LYS
1	A	245	ARG
1	A	279	LEU
1	A	299	LEU
1	A	323	SER
1	A	331	ASN
1	B	28	HIS
1	B	33	GLN
1	B	37	LYS
1	B	40	THR
1	B	48	ASN
1	B	54	TYR
1	B	83	MET
1	B	92	LYS
1	B	98	GLU
1	B	127	ASP
1	B	130	CYS
1	B	148	ASN
1	B	155	LEU
1	B	170	GLU
1	B	180	ASP
1	B	217	ARG
1	B	287	ASP
1	B	315	GLU
1	B	323	SER

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	149	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 monosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 3 ligands modelled in this entry, 3 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	331/334 (99%)	-0.23	2 (0%) 89 88	37, 59, 82, 108	1 (0%)
1	B	330/334 (98%)	0.25	24 (7%) 15 11	38, 75, 114, 139	2 (0%)
All	All	661/668 (98%)	0.01	26 (3%) 39 32	37, 66, 107, 139	3 (0%)

All (26) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	150	LEU	10.8
1	B	152	ASP	8.4
1	B	151	GLY	5.7
1	A	3	LYS	4.5
1	B	180	ASP	3.8
1	B	181	MET	3.8
1	B	78	CYS	3.3
1	B	149	HIS	3.1
1	B	79	ALA	3.1
1	B	176	VAL	2.7
1	B	15	ARG	2.6
1	B	10	SER	2.5
1	B	147	GLY	2.5
1	B	177	HIS	2.5
1	B	184	GLY	2.4
1	B	185	GLN	2.4
1	B	26	ALA	2.4
1	B	47	LYS	2.3
1	B	54	TYR	2.3
1	B	148	ASN	2.2
1	B	175	LEU	2.2
1	B	145	GLN	2.2
1	B	16	ILE	2.1
1	B	146	ILE	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	34	VAL	2.1
1	B	119	ALA	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](#)

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(\AA^2)	Q<0.9
3	MG	A	402	1/1	0.88	0.35	80,80,80,80	0
2	ZN	A	401	1/1	0.98	0.22	49,49,49,49	0
2	ZN	B	401	1/1	0.99	0.26	64,64,64,64	0

6.5 Other polymers [i](#)

There are no such residues in this entry.