



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 05:07 PM UTC

PDB ID : 1CAZ / pdb_00001caz
Title : WILD-TYPE AND E106Q MUTANT CARBONIC ANHYDRASE COM-
PLEXED WITH ACETATE
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Deposited on : 1993-02-26
Resolution : 1.90 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
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

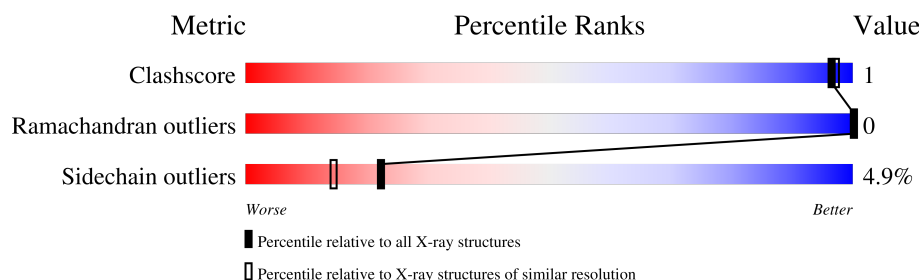
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.90 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))
Clashscore	190562	8410 (1.90-1.90)
Ramachandran outliers	187476	8333 (1.90-1.90)
Sidechain outliers	187428	8333 (1.90-1.90)

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 $\leq 5\%$.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	259	 77% 20%

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	ACY	A	500	-	X	-	-

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 2291 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 CARBONIC ANHYDRASE II.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	258	Total	C	N	O	S	0	4	0
			2075	1330	359	383	3			

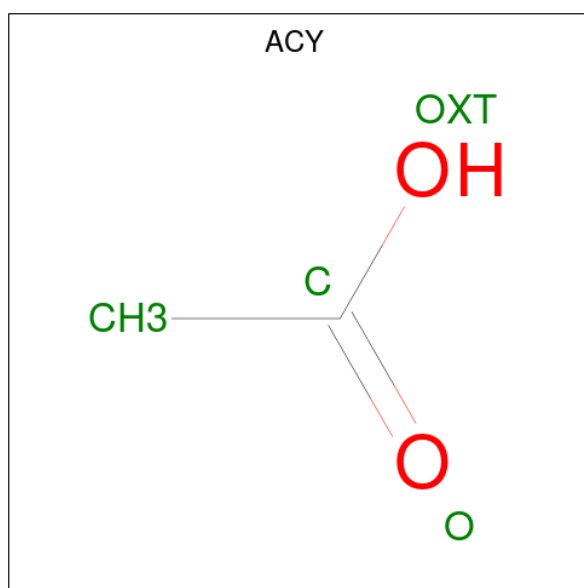
There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	106	GLN	GLU	conflict	UNP P00918

- Molecule 2 is ZINC ION (CCD ID: ZN) (formula: Zn).

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

- Molecule 3 is ACETIC ACID (CCD ID: ACY) (formula: C₂H₄O₂).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			4	2	2		

- Molecule 4 is water.

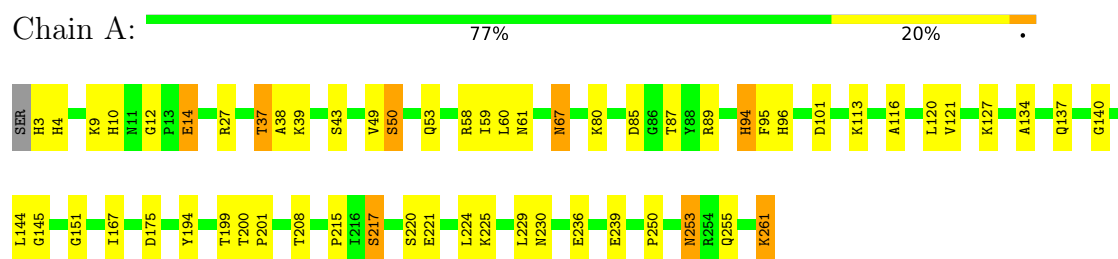
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	210	Total	O	0	0
			210	210		

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. 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. 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.

Note EDS was not executed.

• Molecule 1: CARBONIC ANHYDRASE II



4 Data and refinement statistics

Xtriage (Phenix) and EDS were not executed - this section is therefore incomplete.

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	42.70 Å 41.70 Å 73.00 Å 90.00° 104.60° 90.00°	Depositor
Resolution (Å)	(Not available) – 1.90	Depositor
% Data completeness (in resolution range)	(Not available) ((Not available)-1.90)	Depositor
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	PROFFT	Depositor
R, R_{free}	0.153 , (Not available)	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	2291	wwPDB-VP
Average B, all atoms (Å ²)	12.0	wwPDB-VP

5 Model quality ⓘ

5.1 Standard geometry ⓘ

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

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	1.44	9/2156 (0.4%)	2.01	63/2924 (2.2%)

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	253	ASN	C-O	6.05	1.31	1.24
1	A	200	THR	CA-CB	5.89	1.61	1.54
1	A	94	HIS	CD2-NE2	-5.74	1.31	1.37
1	A	87	THR	CA-CB	5.54	1.61	1.53
1	A	208	THR	C-O	5.48	1.30	1.24
1	A	145	GLY	CA-C	5.40	1.56	1.52
1	A	144	LEU	C-N	-5.31	1.28	1.33
1	A	194	TYR	N-CA	5.31	1.51	1.46
1	A	120	LEU	N-CA	5.17	1.52	1.46

All (63) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	85	ASP	CA-CB-CG	10.62	123.22	112.60
1	A	3	HIS	CA-CB-CG	9.31	123.11	113.80
1	A	96	HIS	CA-CB-CG	-8.49	105.31	113.80
1	A	230	ASN	CA-CB-CG	6.98	119.58	112.60
1	A	10	HIS	CA-CB-CG	-6.91	106.89	113.80
1	A	58	ARG	N-CA-CB	-6.85	100.21	111.66
1	A	49	VAL	CA-C-O	-6.74	112.78	120.66
1	A	144	LEU	CA-C-O	-6.70	112.92	120.43
1	A	37	THR	CB-CA-C	6.55	120.38	109.38
1	A	37	THR	N-CA-CB	-6.54	101.85	110.88
1	A	175	ASP	CB-CA-C	6.42	120.37	109.72
1	A	89	ARG	NE-CZ-NH2	-6.41	113.43	119.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	201	PRO	CB-CA-C	6.39	118.72	110.92
1	A	220	SER	CA-CB-OG	-6.39	98.32	111.10
1	A	217	SER	CB-CA-C	-6.23	99.02	109.48
1	A	120	LEU	CA-C-O	-6.14	114.15	120.54
1	A	27	ARG	CA-CB-CG	6.07	126.23	114.10
1	A	4[A]	HIS	CA-C-N	6.03	129.86	120.75
1	A	4[A]	HIS	C-N-CA	6.03	129.86	120.75
1	A	4[B]	HIS	CA-C-N	6.03	129.86	120.75
1	A	4[B]	HIS	C-N-CA	6.03	129.86	120.75
1	A	250	PRO	CB-CA-C	-6.00	103.72	111.46
1	A	87	THR	CA-CB-OG1	-6.00	100.61	109.60
1	A	89	ARG	O-C-N	5.96	130.01	123.22
1	A	94	HIS	CA-CB-CG	-5.96	107.84	113.80
1	A	236	GLU	CA-CB-CG	5.93	125.96	114.10
1	A	221	GLU	CG-CD-OE1	5.78	131.70	118.40
1	A	58	ARG	O-C-N	5.75	129.81	123.25
1	A	60	LEU	N-CA-CB	-5.74	100.85	111.13
1	A	137	GLN	O-C-N	5.66	127.06	121.57
1	A	14[A]	GLU	CB-CG-CD	5.61	122.14	112.60
1	A	14[B]	GLU	CB-CG-CD	5.61	122.14	112.60
1	A	38	ALA	N-CA-C	-5.59	102.01	110.28
1	A	95	PHE	O-C-N	5.56	130.56	122.94
1	A	230	ASN	CB-CG-ND2	5.52	124.68	116.40
1	A	67	ASN	CA-CB-CG	5.52	118.12	112.60
1	A	113	LYS	CA-C-O	-5.52	114.77	121.28
1	A	225	LYS	CA-CB-CG	-5.51	103.07	114.10
1	A	224	LEU	CA-C-O	-5.48	114.71	120.63
1	A	101	ASP	CA-CB-CG	5.46	118.06	112.60
1	A	253	ASN	CA-C-O	-5.45	112.72	120.51
1	A	89	ARG	CB-CA-C	-5.43	100.89	109.80
1	A	85	ASP	N-CA-CB	5.43	119.12	110.65
1	A	199	THR	OG1-CB-CG2	5.42	120.14	109.30
1	A	261	LYS	CA-CB-CG	5.37	124.84	114.10
1	A	121	VAL	O-C-N	5.35	128.96	123.03
1	A	127	LYS	CA-CB-CG	5.30	124.70	114.10
1	A	50	SER	N-CA-CB	5.29	119.41	110.46
1	A	215	PRO	CA-C-O	-5.17	115.71	122.12
1	A	53	GLN	CG-CD-NE2	-5.14	108.68	116.40
1	A	61	ASN	N-CA-C	-5.13	99.42	108.20
1	A	53	GLN	OE1-CD-NE2	5.12	127.72	122.60
1	A	12	GLY	O-C-N	5.12	126.89	121.77
1	A	208	THR	N-CA-C	-5.09	100.43	108.73

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	116	ALA	O-C-N	5.08	128.60	123.26
1	A	229	LEU	O-C-N	5.08	128.90	122.65
1	A	37	THR	CA-CB-CG2	5.07	119.11	110.50
1	A	95	PHE	CA-C-O	-5.06	115.71	121.58
1	A	151	GLY	N-CA-C	-5.06	101.19	113.18
1	A	236	GLU	CG-CD-OE1	5.02	129.95	118.40
1	A	217	SER	CA-CB-OG	-5.01	101.08	111.10
1	A	89	ARG	NE-CZ-NH1	5.00	126.50	121.50
1	A	261	LYS	N-CA-CB	5.00	119.00	110.50

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	2075	0	2021	4	0
2	A	2	0	0	0	0
3	A	4	0	3	0	0
4	A	210	0	0	1	0
All	All	2291	0	2024	4	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 1.

All (4) 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:59:ILE:HG12	1:A:167:ILE:HD13	1.95	0.48
1:A:134:ALA:O	1:A:140:GLY:HA3	2.16	0.45
1:A:67:ASN:HD22	1:A:94:HIS:HB3	1.84	0.43
1:A:80:LYS:HG2	4:A:416:HOH:O	2.20	0.42

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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	260/259 (100%)	249 (96%)	11 (4%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains

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	227/224 (101%)	215 (95%)	12 (5%)	20	12

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

Mol	Chain	Res	Type
1	A	9	LYS
1	A	14[A]	GLU
1	A	14[B]	GLU
1	A	37	THR
1	A	39	LYS
1	A	43	SER
1	A	50	SER
1	A	217	SER
1	A	239	GLU
1	A	253	ASN
1	A	255	GLN
1	A	261	LYS

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

Mol	Chain	Res	Type
1	A	67	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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 3 ligands modelled in this entry, 2 are monoatomic - leaving 1 for Mogul analysis.

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	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
3	ACY	A	500	2	3,3,3	1.70	1 (33%)	3,3,3	5.24	3 (100%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	500	ACY	OXT-C	-2.10	1.21	1.30

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
3	A	500	ACY	OXT-C-CH3	7.26	145.49	115.05
3	A	500	ACY	OXT-C-O	-3.95	107.36	122.03
3	A	500	ACY	O-C-CH3	-3.76	107.12	122.53

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

EDS was not executed - this section is therefore empty.

6.2 Non-standard residues in protein, DNA, RNA chains

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates

EDS was not executed - this section is therefore empty.

6.4 Ligands

EDS was not executed - this section is therefore empty.

6.5 Other polymers

EDS was not executed - this section is therefore empty.