



# wwPDB X-ray Structure Validation Summary Report ⓘ

Apr 13, 2026 – 02:58 PM JST

PDB ID : 9X6A / pdb\_00009x6a  
Title : carboxylesterase EstC  
Authors : Chen, Q.; Li, B.  
Deposited on : 2025-10-15  
Resolution : 2.06 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](mailto: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>.

---

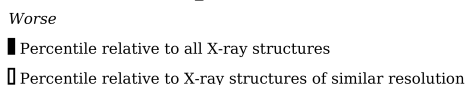
The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

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

**i**

## X-RAY DIFFRACTION

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 <sub>free</sub>	180053	3774 (2.08-2.04)
Clashscore	190562	3883 (2.08-2.04)
Ramachandran outliers	187476	3860 (2.08-2.04)
Sidechain outliers	187428	3860 (2.08-2.04)
RSRZ outliers	180081	3775 (2.08-2.04)

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  $\geq 3$ , 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	301	<div> <div>4%</div> <div>88%</div> <div>7%</div> <div></div> </div>
1	B	301	<div> <div>4%</div> <div>88%</div> <div>7%</div> <div></div> </div>
1	C	301	<div> <div>5%</div> <div>84%</div> <div>11%</div> <div>5%</div> </div>
1	D	301	<div> <div>7%</div> <div>81%</div> <div>12%</div> <div>5%</div> </div>
1	E	301	<div> <div>4%</div> <div>82%</div> <div>11%</div> <div>6%</div> </div>

## 2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 11186 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 carboxylesterase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	288	Total	C	N	O	S	0	0	0
			2119	1337	385	395	2			
1	B	289	Total	C	N	O	S	0	0	0
			2123	1339	386	396	2			
1	C	285	Total	C	N	O	S	0	0	0
			2103	1327	382	392	2			
1	D	285	Total	C	N	O	S	0	0	0
			2100	1325	382	391	2			
1	E	283	Total	C	N	O	S	0	0	0
			2090	1320	380	388	2			

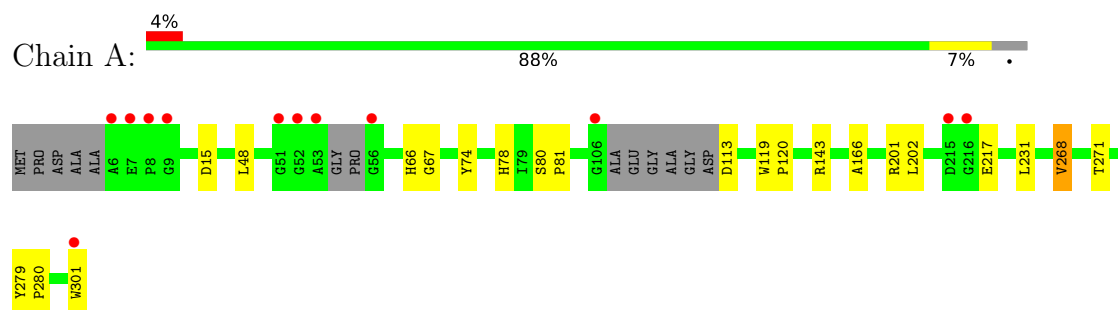
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	165	Total	O	0	0
			165	165		
2	B	141	Total	O	0	0
			141	141		
2	C	142	Total	O	0	0
			142	142		
2	D	98	Total	O	0	0
			98	98		
2	E	105	Total	O	0	0
			105	105		

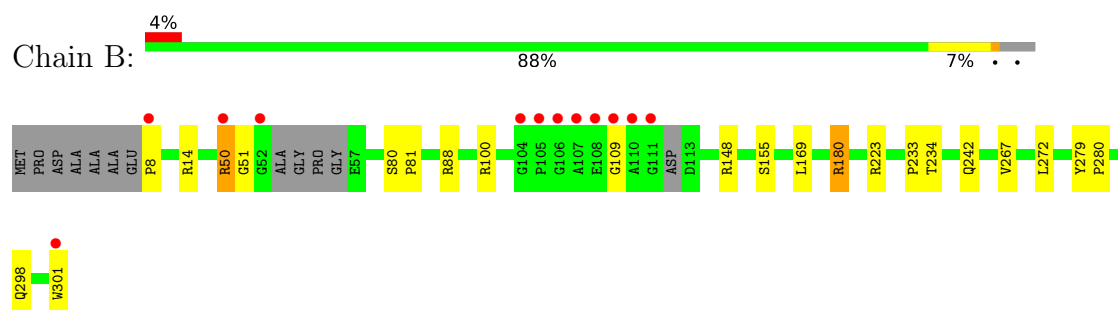
### 3 Residue-property plots [i](#)

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.

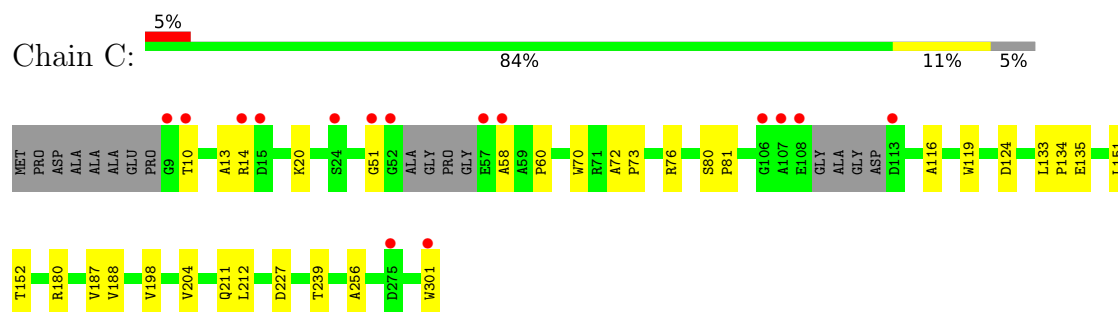
- Molecule 1: carboxylesterase



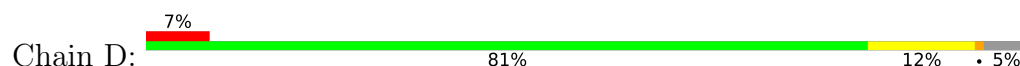
- Molecule 1: carboxylesterase

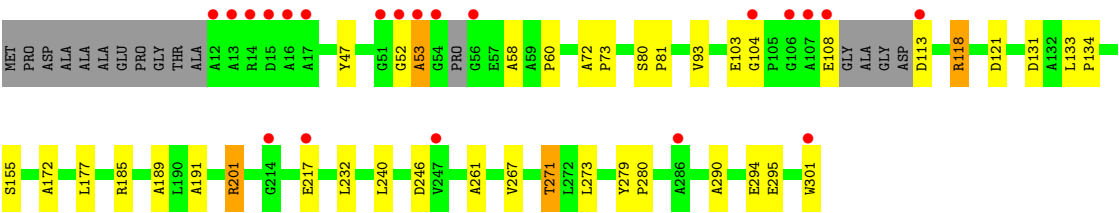


- Molecule 1: carboxylesterase

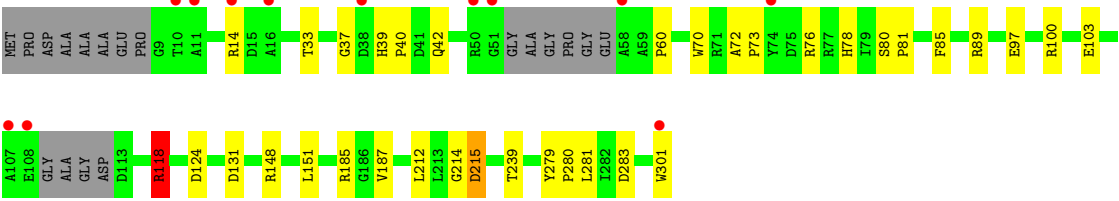
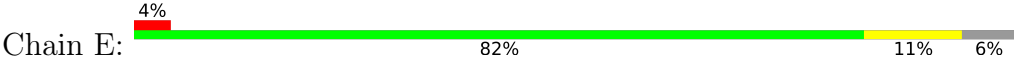


- Molecule 1: carboxylesterase





● Molecule 1: carboxylesterase



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	47.48Å 108.27Å 184.32Å 90.00° 90.70° 90.00°	Depositor
Resolution (Å)	27.37 – 2.06 27.37 – 2.06	Depositor EDS
% Data completeness (in resolution range)	99.3 (27.37-2.06) 99.3 (27.37-2.06)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.51 (at 2.06Å)	Xtriage
Refinement program	REFMAC 5.8.0430	Depositor
R, $R_{free}$	0.212 , 0.248 0.218 , 0.253	Depositor DCC
$R_{free}$ test set	5706 reflections (4.99%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	23.1	Xtriage
Anisotropy	0.030	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.38 , 38.0	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.43$ , $\langle L^2 \rangle = 0.26$	Xtriage
Estimated twinning fraction	0.055 for h,-k,-l	Xtriage
$F_o, F_c$ correlation	0.92	EDS
Total number of atoms	11186	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	26.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.33% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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	0.68	0/2173	1.12	5/2977 (0.2%)
1	B	0.69	0/2177	1.11	2/2981 (0.1%)
1	C	0.64	0/2156	1.08	3/2953 (0.1%)
1	D	0.62	0/2153	1.11	4/2948 (0.1%)
1	E	0.63	0/2143	1.07	6/2936 (0.2%)
All	All	0.65	0/10802	1.10	20/14795 (0.1%)

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	A	0	1
1	B	0	3
1	D	0	1
1	E	0	2
All	All	0	7

There are no bond length outliers.

The worst 5 of 20 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	268	VAL	N-CA-CB	-7.58	101.90	111.46
1	E	124	ASP	CA-CB-CG	6.83	119.44	112.60
1	B	298	GLN	CB-CA-C	-6.54	99.73	110.85
1	E	131	ASP	CA-CB-CG	6.15	118.75	112.60
1	D	131	ASP	CA-CB-CG	6.04	118.64	112.60

There are no chirality outliers.

5 of 7 planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	201	ARG	Sidechain
1	B	100	ARG	Sidechain
1	B	14	ARG	Sidechain
1	B	223	ARG	Sidechain
1	D	113	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	2119	0	2057	7	0
1	B	2123	0	2061	8	0
1	C	2103	0	2042	14	0
1	D	2100	0	2038	18	0
1	E	2090	0	2033	17	0
2	A	165	0	0	1	0
2	B	141	0	0	0	0
2	C	142	0	0	1	0
2	D	98	0	0	0	0
2	E	105	0	0	1	0
All	All	11186	0	10231	64	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.

The worst 5 of 64 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:E:103:GLU:HB2	1:E:118:ARG:HD2	1.61	0.81
1:B:180:ARG:HH11	1:B:180:ARG:HB2	1.54	0.73
1:D:201:ARG:HH11	1:D:201:ARG:HG3	1.53	0.72
1:D:271:THR:HG21	1:D:295:GLU:OE2	1.94	0.66
1:E:281:LEU:HD11	2:E:441:HOH:O	2.01	0.60

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	282/301 (94%)	275 (98%)	7 (2%)	0	100	100
1	B	283/301 (94%)	272 (96%)	9 (3%)	2 (1%)	18	10
1	C	279/301 (93%)	269 (96%)	8 (3%)	2 (1%)	18	10
1	D	279/301 (93%)	265 (95%)	12 (4%)	2 (1%)	18	10
1	E	277/301 (92%)	269 (97%)	7 (2%)	1 (0%)	30	23
All	All	1400/1505 (93%)	1350 (96%)	43 (3%)	7 (0%)	24	17

5 of 7 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	53	ALA
1	B	51	GLY
1	D	58	ALA
1	E	215	ASP
1	B	109	GLY

### 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	200/206 (97%)	196 (98%)	4 (2%)	48	47
1	B	200/206 (97%)	195 (98%)	5 (2%)	42	38
1	C	199/206 (97%)	195 (98%)	4 (2%)	48	47
1	D	198/206 (96%)	192 (97%)	6 (3%)	36	32

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	E	198/206 (96%)	197 (100%)	1 (0%)	81	85
All	All	995/1030 (97%)	975 (98%)	20 (2%)	48	47

5 of 20 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	D	201	ARG
1	D	271	THR
1	E	78	HIS
1	D	273	LEU
1	B	155	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	139	GLN
1	D	139	GLN
1	D	251	GLN
1	E	139	GLN
1	E	298	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](#)

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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	288/301 (95%)	0.03	12 (4%)	40 41	12, 19, 42, 70	0
1	B	289/301 (96%)	0.00	12 (4%)	40 41	10, 18, 39, 80	0
1	C	285/301 (94%)	0.25	15 (5%)	32 32	13, 23, 47, 75	0
1	D	285/301 (94%)	0.55	21 (7%)	20 20	14, 28, 55, 80	0
1	E	283/301 (94%)	0.38	12 (4%)	40 41	16, 26, 49, 76	0
All	All	1430/1505 (95%)	0.24	72 (5%)	34 34	10, 23, 49, 80	0

The worst 5 of 72 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	107	ALA	6.8
1	D	13	ALA	5.5
1	A	53	ALA	5.1
1	E	107	ALA	5.0
1	C	51	GLY	4.9

### 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](#)

There are no ligands in this entry.

## 6.5 Other polymers [i](#)

There are no such residues in this entry.