



# wwPDB X-ray Structure Validation Summary Report ⓘ

Feb 6, 2024 – 01:49 AM EST

PDB ID : 1Z7L  
Title : Crystal structure of fragment of mouse ubiquitin-activating enzyme  
Authors : Szczepanowski, R.H.; Filipek, R.; Bochtler, M.  
Deposited on : 2005-03-25  
Resolution : 2.80 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>.

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

MolProbity : 4.02b-467  
Mogul : 1.8.5 (274361), CSD as541be (2020)  
Xtriage (Phenix) : 1.13  
EDS : 2.36  
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) : 2.36

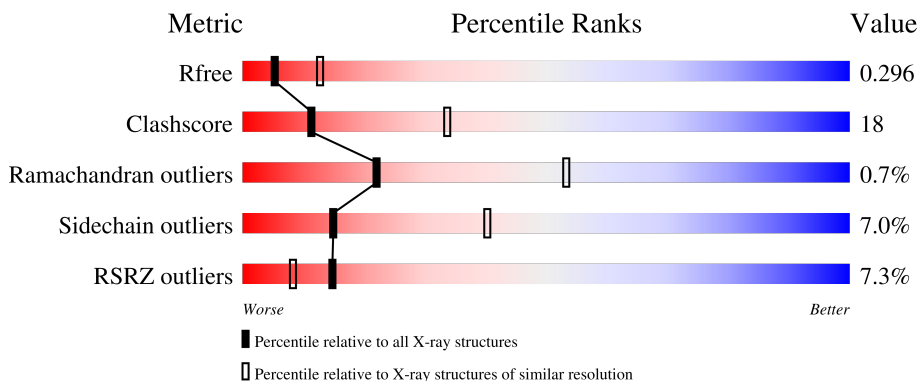
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.80 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	3140 (2.80-2.80)
Clashscore	141614	3569 (2.80-2.80)
Ramachandran outliers	138981	3498 (2.80-2.80)
Sidechain outliers	138945	3500 (2.80-2.80)
RSRZ outliers	127900	3078 (2.80-2.80)

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	276	 7% 62% 28% 8%
1	B	276	 3% 63% 27% 8%
1	C	276	 10% 62% 28% 8%

## 2 Entry composition i

There are 2 unique types of molecules in this entry. The entry contains 6144 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 Ubiquitin-activating enzyme E1 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	255	2030	1287	353	382	8	0	0	0
1	B	255	2030	1287	353	382	8	0	0	0
1	C	255	2030	1287	353	382	8	0	0	0

There are 30 discrepancies between the modelled and reference sequences:

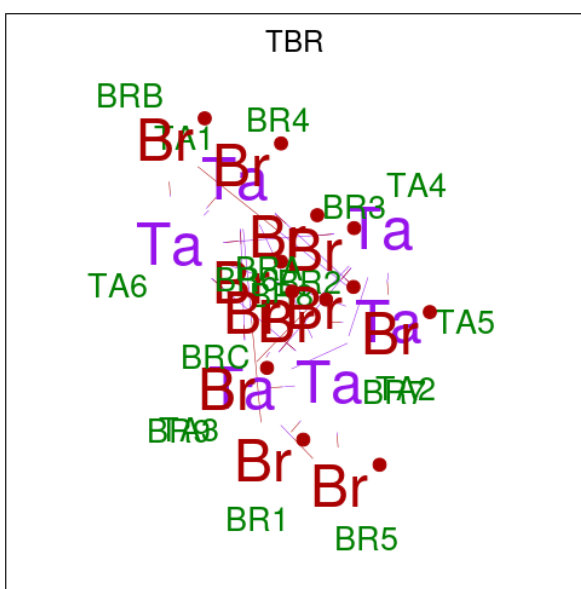
Chain	Residue	Modelled	Actual	Comment	Reference
A	616	MET	-	expression tag	UNP Q02053
A	617	GLY	-	expression tag	UNP Q02053
A	618	HIS	-	expression tag	UNP Q02053
A	619	HIS	-	expression tag	UNP Q02053
A	620	HIS	-	expression tag	UNP Q02053
A	621	HIS	-	expression tag	UNP Q02053
A	622	HIS	-	expression tag	UNP Q02053
A	623	HIS	-	expression tag	UNP Q02053
A	624	GLU	-	expression tag	UNP Q02053
A	625	PHE	-	expression tag	UNP Q02053
B	616	MET	-	expression tag	UNP Q02053
B	617	GLY	-	expression tag	UNP Q02053
B	618	HIS	-	expression tag	UNP Q02053
B	619	HIS	-	expression tag	UNP Q02053
B	620	HIS	-	expression tag	UNP Q02053
B	621	HIS	-	expression tag	UNP Q02053
B	622	HIS	-	expression tag	UNP Q02053
B	623	HIS	-	expression tag	UNP Q02053
B	624	GLU	-	expression tag	UNP Q02053
B	625	PHE	-	expression tag	UNP Q02053
C	616	MET	-	expression tag	UNP Q02053
C	617	GLY	-	expression tag	UNP Q02053
C	618	HIS	-	expression tag	UNP Q02053

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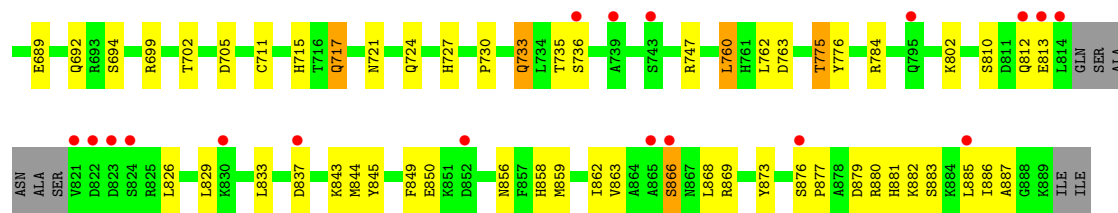
Chain	Residue	Modelled	Actual	Comment	Reference
C	619	HIS	-	expression tag	UNP Q02053
C	620	HIS	-	expression tag	UNP Q02053
C	621	HIS	-	expression tag	UNP Q02053
C	622	HIS	-	expression tag	UNP Q02053
C	623	HIS	-	expression tag	UNP Q02053
C	624	GLU	-	expression tag	UNP Q02053
C	625	PHE	-	expression tag	UNP Q02053

- Molecule 2 is HEXATANTALUM DODECABROMIDE (three-letter code: TBR) (formula:  $\text{Br}_{12}\text{Ta}_6$ ).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	Br	Ta		
2	A	1	18	12	6	0	0
2	B	1	18	12	6	0	0
2	C	1	18	12	6	0	0





## 4 Data and refinement statistics

Property	Value	Source
Space group	H 3 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	216.12Å 216.12Å 196.09Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	10.00 – 2.80 19.99 – 2.80	Depositor EDS
% Data completeness (in resolution range)	99.0 (10.00-2.80) 99.0 (19.99-2.80)	Depositor EDS
$R_{merge}$	0.07	Depositor
$R_{sym}$	0.07	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	5.06 (at 2.79Å)	Xtrriage
Refinement program	CNS 1.0	Depositor
R, $R_{free}$	0.277 , 0.308 0.267 , 0.296	Depositor DCC
$R_{free}$ test set	2109 reflections (4.93%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	73.9	Xtrriage
Anisotropy	0.012	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.32 , 57.8	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.33$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
$F_o, F_c$ correlation	0.91	EDS
Total number of atoms	6144	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	76.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.68% 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

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

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	0.41	0/2085	0.65	0/2841
1	B	0.42	0/2085	0.66	0/2841
1	C	0.42	0/2085	0.65	0/2841
All	All	0.42	0/6255	0.65	0/8523

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

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	2030	0	1953	71	0
1	B	2030	0	1953	70	0
1	C	2030	0	1953	72	0
2	A	18	0	0	3	0
2	B	18	0	0	3	0
2	C	18	0	0	2	0
All	All	6144	0	5859	213	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 18.



The worst 5 of 213 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:717:GLN:HE21	1:A:717:GLN:HA	1.15	1.09
1:B:717:GLN:HA	1:B:717:GLN:HE21	1.20	1.05
1:C:717:GLN:HA	1:C:717:GLN:HE21	1.15	1.03
1:C:649:ARG:HH12	1:C:653:GLU:HG2	1.24	1.01
1:A:649:ARG:HH12	1:A:653:GLU:HG2	1.27	0.98

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	251/276 (91%)	232 (92%)	18 (7%)	1 (0%)	34	66
1	B	251/276 (91%)	235 (94%)	14 (6%)	2 (1%)	19	49
1	C	251/276 (91%)	233 (93%)	16 (6%)	2 (1%)	19	49
All	All	753/828 (91%)	700 (93%)	48 (6%)	5 (1%)	22	53

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	681	GLY
1	B	681	GLY
1	C	681	GLY
1	C	680	ALA
1	B	680	ALA

### 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	223/243 (92%)	207 (93%)	16 (7%)	14	38
1	B	223/243 (92%)	209 (94%)	14 (6%)	18	46
1	C	223/243 (92%)	206 (92%)	17 (8%)	13	36
All	All	669/729 (92%)	622 (93%)	47 (7%)	15	40

5 of 47 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	813	GLU
1	C	684	PRO
1	B	837	ASP
1	C	646	GLN
1	C	699	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 11 such sidechains are listed below:

Mol	Chain	Res	Type
1	B	717	GLN
1	C	643	HIS
1	C	717	GLN
1	C	715	HIS
1	B	643	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](#)

3 ligands are modelled in this entry.

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
2	TBR	C	1001	-	0,36,36	-	-	-		
2	TBR	A	999	-	0,36,36	-	-	-		
2	TBR	B	1000	-	0,36,36	-	-	-		

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.

3 monomers are involved in 8 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	1001	TBR	2	0
2	A	999	TBR	3	0
2	B	1000	TBR	3	0

## 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	255/276 (92%)	0.32	20 (7%) 13 7	23, 70, 112, 161	0
1	B	255/276 (92%)	0.14	9 (3%) 44 34	23, 63, 103, 162	0
1	C	255/276 (92%)	0.62	27 (10%) 6 3	23, 81, 117, 162	0
All	All	765/828 (92%)	0.36	56 (7%) 15 8	23, 73, 113, 162	0

The worst 5 of 56 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	821	VAL	7.2
1	B	821	VAL	6.5
1	C	629	ILE	4.3
1	C	812	GLN	4.1
1	B	885	LEU	4.0

### 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, 95<sup>th</sup> 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	TBR	A	999	18/18	0.84	0.25	83,86,87,88	18
2	TBR	B	1000	18/18	0.85	0.27	86,87,88,88	18
2	TBR	C	1001	18/18	0.89	0.17	86,87,88,89	18

## 6.5 Other polymers [i](#)

There are no such residues in this entry.