



# Full wwPDB X-ray Structure Validation Report ⓘ

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PDB ID : 7WG2  
Title : EVAA-KlAtel  
Authors : Kim, M.K.; Kim, B.H.; Oh, S.-J.; Song, H.K.  
Deposited on : 2021-12-28  
Resolution : 1.76 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](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  
Xtriage (Phenix) : 1.13  
EDS : 2.30  
buster-report : 1.1.7 (2018)  
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) : 2.30

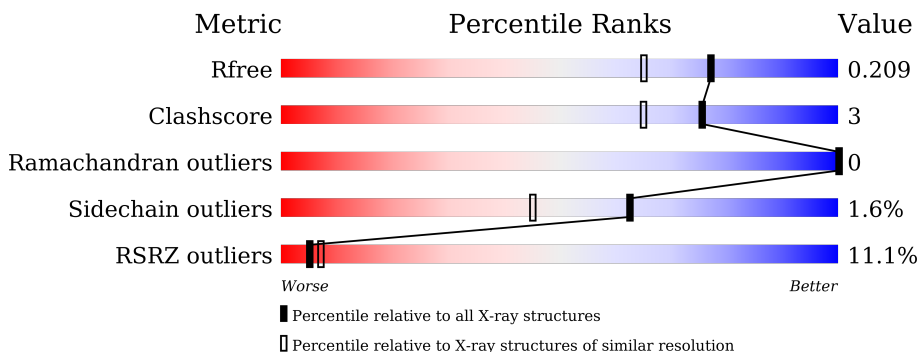
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 1.76 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	2340 (1.76-1.76)
Clashscore	141614	2466 (1.76-1.76)
Ramachandran outliers	138981	2437 (1.76-1.76)
Sidechain outliers	138945	2437 (1.76-1.76)
RSRZ outliers	127900	2298 (1.76-1.76)

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	507	

## 2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 8168 atoms, of which 3880 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 Arginyltransferase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace	
			Total	C	H	N	O				S
1	A	487	7828	2533	3880	643	755	17	0	0	0

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-3	GLU	-	expression tag	UNP Q6CXX6
A	-2	VAL	-	expression tag	UNP Q6CXX6
A	-1	ALA	-	expression tag	UNP Q6CXX6
A	0	ALA	-	expression tag	UNP Q6CXX6

- Molecule 2 is ZINC ION (three-letter code: ZN) (formula: Zn) (labeled as "Ligand of Interest" by depositor).

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

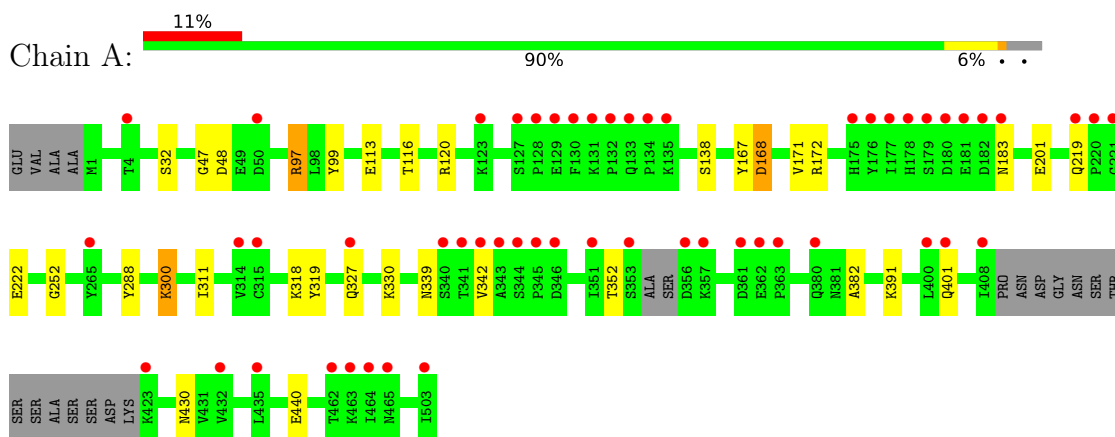
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	339	Total	O	0	0
			339	339		

### 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: Arginyltransferase



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 43 21 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	86.60Å 86.60Å 161.11Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	29.49 – 1.76 29.49 – 1.76	Depositor EDS
% Data completeness (in resolution range)	100.0 (29.49-1.76) 100.0 (29.49-1.76)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.72 (at 1.76Å)	Xtrriage
Refinement program	PHENIX 1.19.2_4158	Depositor
R, $R_{free}$	0.179 , 0.212 0.177 , 0.209	Depositor DCC
$R_{free}$ test set	3073 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	25.9	Xtrriage
Anisotropy	0.303	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.39 , 42.3	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.34$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
$F_o, F_c$ correlation	0.96	EDS
Total number of atoms	8168	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	36.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.32% 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 [i](#)

### 5.1 Standard geometry [i](#)

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

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	0.51	0/4043	0.67	0/5473

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

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	97	ARG	Sidechain

### 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	3948	3880	3882	24	0
2	A	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	A	339	0	0	6	0
All	All	4288	3880	3882	24	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.

All (24) 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:300:LYS:HE2	1:A:300:LYS:H	1.46	0.81
1:A:327:GLN:OE1	1:A:330:LYS:NZ	2.12	0.80
1:A:339:ASN:O	1:A:342:VAL:HG12	1.93	0.68
1:A:318:LYS:HD3	1:A:319:TYR:H	1.61	0.66
1:A:183:ASN:N	3:A:701:HOH:O	2.13	0.64
1:A:99:TYR:HB3	1:A:311:ILE:HG23	1.82	0.60
1:A:168:ASP:OD1	1:A:172:ARG:NH2	2.35	0.57
1:A:219:GLN:N	1:A:222:GLU:OE2	2.28	0.57
1:A:300:LYS:H	1:A:300:LYS:CE	2.20	0.54
1:A:318:LYS:HD3	1:A:319:TYR:N	2.23	0.53
1:A:318:LYS:CD	1:A:319:TYR:N	2.79	0.46
1:A:252:GLY:HA3	1:A:288:TYR:O	2.16	0.45
1:A:201:GLU:OE2	3:A:702:HOH:O	2.21	0.45
1:A:401:GLN:H	1:A:401:GLN:CD	2.21	0.44
1:A:168:ASP:O	1:A:171:VAL:HG22	2.19	0.43
1:A:391:LYS:HB3	3:A:836:HOH:O	2.18	0.43
1:A:183:ASN:HB2	3:A:701:HOH:O	2.19	0.42
1:A:352:THR:N	1:A:440:GLU:OE2	2.43	0.41
1:A:47:GLY:O	3:A:703:HOH:O	2.22	0.41
1:A:167:TYR:O	1:A:171:VAL:HG13	2.19	0.41
1:A:219:GLN:CB	1:A:222:GLU:OE2	2.69	0.41
1:A:382:ALA:HB1	1:A:430:ASN:HB3	2.02	0.41
1:A:352:THR:HB	3:A:988:HOH:O	2.21	0.40
1:A:116:THR:O	1:A:120:ARG:HG3	2.20	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	481/507 (95%)	472 (98%)	9 (2%)	0	100	100

There are no Ramachandran outliers to report.

### 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	446/461 (97%)	439 (98%)	7 (2%)	62	45

All (7) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	32	SER
1	A	48	ASP
1	A	97	ARG
1	A	113	GLU
1	A	138	SER
1	A	168	ASP
1	A	300	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.



### 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 1 ligands modelled in this entry, 1 is 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, 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	487/507 (96%)	0.52	54 (11%) <b>5</b> <b>7</b>	16, 28, 60, 98	0

All (54) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	464	ILE	8.1
1	A	408	ILE	6.1
1	A	363	PRO	5.5
1	A	181	GLU	5.4
1	A	128	PRO	5.4
1	A	343	ALA	5.2
1	A	180	ASP	5.2
1	A	463	LYS	5.1
1	A	182	ASP	4.9
1	A	131	LYS	4.6
1	A	130	PHE	4.6
1	A	342	VAL	4.4
1	A	134	PRO	4.4
1	A	346	ASP	3.9
1	A	4	THR	3.8
1	A	503	ILE	3.7
1	A	129	GLU	3.6
1	A	340	SER	3.6
1	A	341	THR	3.6
1	A	133	GLN	3.5
1	A	356	ASP	3.4
1	A	179	SER	3.4
1	A	221	GLY	3.3
1	A	462	THR	3.2
1	A	178	HIS	3.1
1	A	127	SER	3.0
1	A	361	ASP	2.9

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Mol	Chain	Res	Type	RSRZ
1	A	183	ASN	2.7
1	A	219	GLN	2.7
1	A	380	GLN	2.7
1	A	265	TYR	2.6
1	A	435	LEU	2.5
1	A	400	LEU	2.5
1	A	176	TYR	2.5
1	A	344	SER	2.5
1	A	123	LYS	2.4
1	A	132	PRO	2.4
1	A	353	SER	2.3
1	A	465	ASN	2.3
1	A	135	LYS	2.2
1	A	50	ASP	2.2
1	A	327	GLN	2.2
1	A	220	PRO	2.2
1	A	401	GLN	2.2
1	A	432	VAL	2.1
1	A	314	VAL	2.1
1	A	175	HIS	2.1
1	A	351	ILE	2.1
1	A	423	LYS	2.1
1	A	315	CYS	2.1
1	A	177	ILE	2.1
1	A	362	GLU	2.0
1	A	357	LYS	2.0
1	A	345	PRO	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.

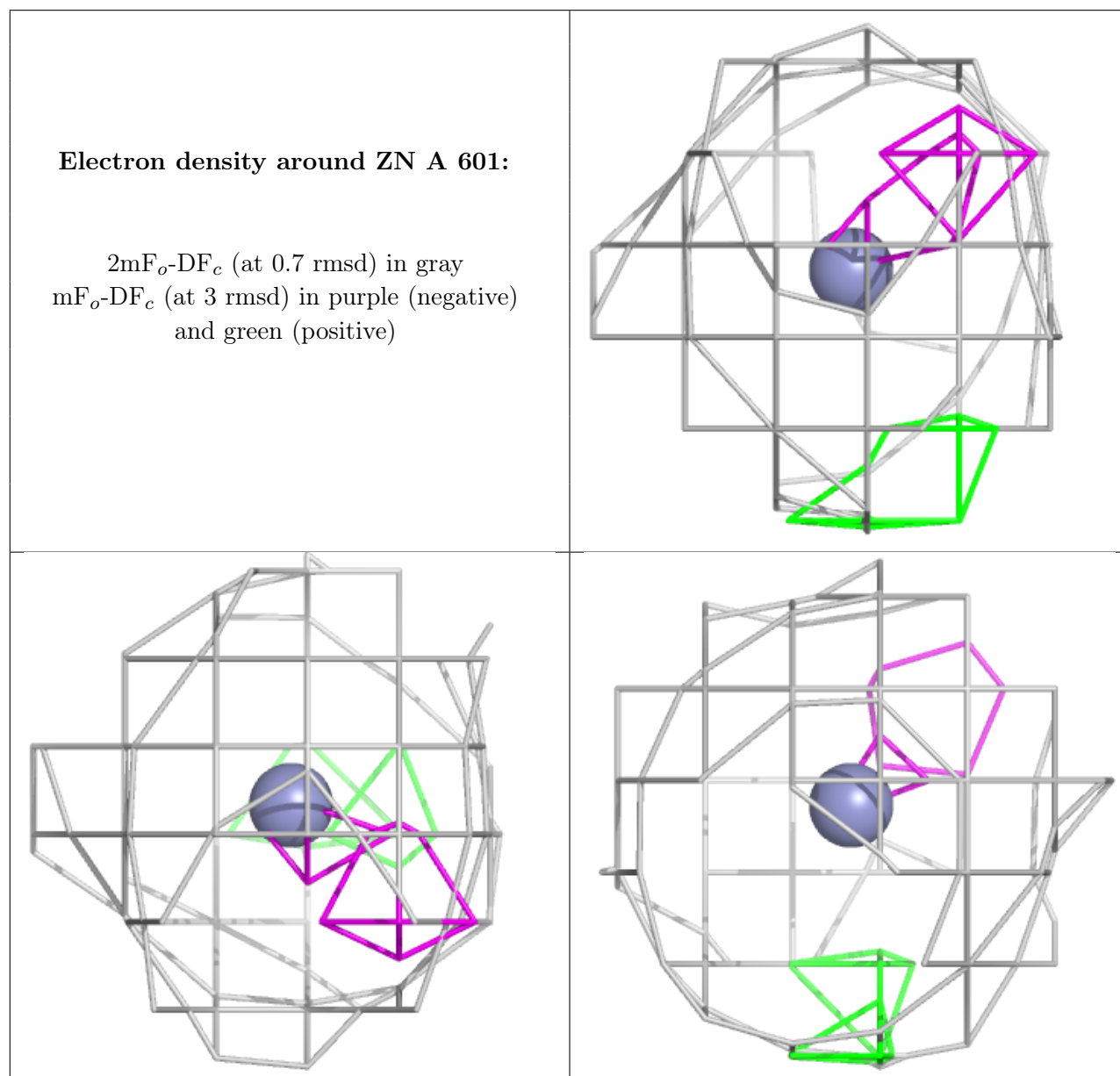
## 6.4 Ligands [i](#)

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, 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(Å <sup>2</sup> )	Q<0.9
2	ZN	A	601	1/1	1.00	0.07	21,21,21,21	0

The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.



## 6.5 Other polymers [i](#)

There are no such residues in this entry.