



# Full wwPDB X-ray Structure Validation Report ⓘ

May 21, 2020 – 06:09 am BST

PDB ID : 5ISW  
Title : Structure of the apo PCP-E didomain of the gramicidin S synthetase A  
Authors : Chen, W.-H.; Li, K.; Bruner, S.D.  
Deposited on : 2016-03-15  
Resolution : 1.75 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.

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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 : 2.11  
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) : 2.11

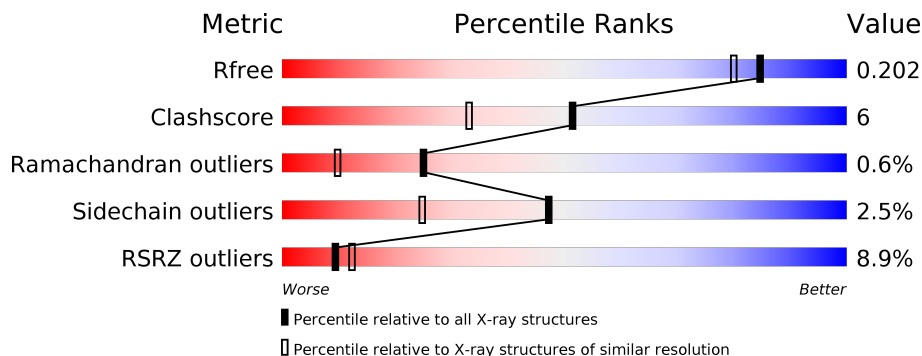
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 1.75 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 on 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	573	

## 2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 4955 atoms, of which 8 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 Gramicidin S synthase 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	539	4471	2867	749	842	13	0	2	0

There are 13 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	537	MET	-	initiating methionine	UNP P0C062
A	1009	GLY	SER	conflict	UNP P0C062
A	1099	ASP	-	expression tag	UNP P0C062
A	1100	ASP	-	expression tag	UNP P0C062
A	1101	ASP	-	expression tag	UNP P0C062
A	1102	ASP	-	expression tag	UNP P0C062
A	1103	LYS	-	expression tag	UNP P0C062
A	1104	HIS	-	expression tag	UNP P0C062
A	1105	HIS	-	expression tag	UNP P0C062
A	1106	HIS	-	expression tag	UNP P0C062
A	1107	HIS	-	expression tag	UNP P0C062
A	1108	HIS	-	expression tag	UNP P0C062
A	1109	HIS	-	expression tag	UNP P0C062

- Molecule 2 is GLYCEROL (three-letter code: GOL) (formula: C<sub>3</sub>H<sub>8</sub>O<sub>3</sub>).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	C	H	O		
2	A	1	14	3	8	3	0	0

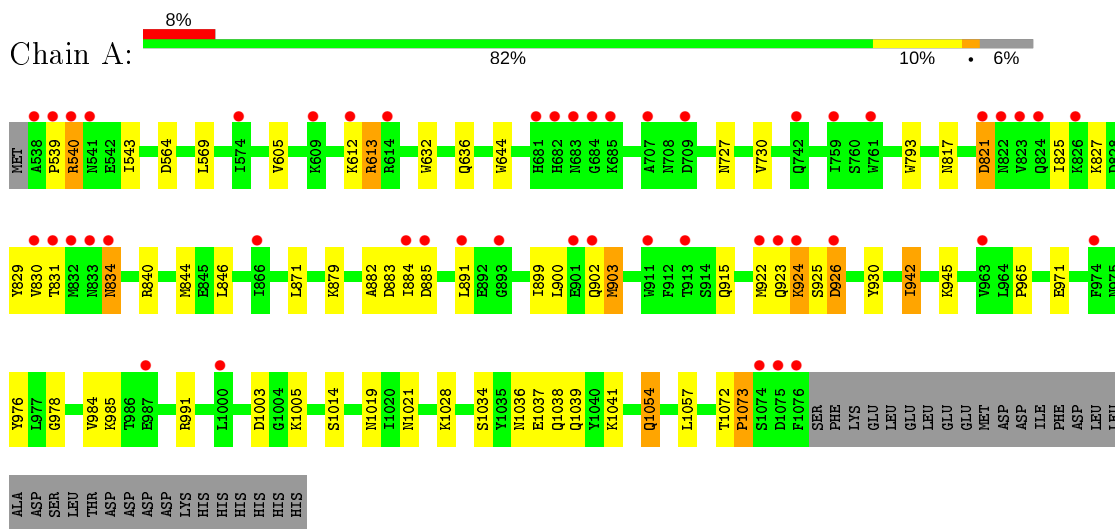
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	470	Total	O	0	0
			470	470		

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

These plots are drawn for all protein, RNA and DNA 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: Gramicidin S synthase 1



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	53.77Å 57.64Å 61.38Å 65.72° 65.47° 84.88°	Depositor
Resolution (Å)	34.00 – 1.75 34.00 – 1.75	Depositor EDS
% Data completeness (in resolution range)	97.2 (34.00-1.75) 93.8 (34.00-1.75)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.76 (at 1.75Å)	Xtrriage
Refinement program	PHENIX 1.9_1692	Depositor
R, $R_{free}$	0.169 , 0.201 0.173 , 0.202	Depositor DCC
$R_{free}$ test set	1990 reflections (3.34%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	20.1	Xtrriage
Anisotropy	0.134	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.40 , 48.1	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.48$ , $\langle L^2 \rangle = 0.31$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
$F_o, F_c$ correlation	0.96	EDS
Total number of atoms	4955	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	26.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.90% 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: GOL

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	0.58	0/4577	0.61	0/6192

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 [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	4471	0	4388	50	0
2	A	6	8	8	0	0
3	A	470	0	0	6	0
All	All	4947	8	4396	50	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 (50) 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:827:LYS:HG2	1:A:1039:GLN:HG2	1.69	0.75

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:831:THR:HG21	1:A:1038:GLN:O	1.92	0.70
1:A:612:LYS:O	1:A:613:ARG:HB2	1.96	0.66
1:A:827:LYS:HG2	1:A:1039:GLN:CG	2.27	0.65
1:A:825:ILE:HG22	1:A:884:ILE:HD11	1.80	0.64
1:A:924:LYS:HG3	1:A:930:TYR:CD2	2.33	0.64
1:A:834:ASN:HB3	1:A:1036:ASN:ND2	2.14	0.62
1:A:882:ALA:HB3	1:A:884:ILE:HG13	1.80	0.62
1:A:925:SER:O	1:A:926:ASP:HB2	2.01	0.59
1:A:985:LYS:HE2	1:A:991:ARG:HH12	1.70	0.56
1:A:902:GLN:O	1:A:903:MET:HE2	2.07	0.54
1:A:840:ARG:HG3	1:A:1037:GLU:HG3	1.89	0.54
1:A:1072:THR:HB	1:A:1073:PRO:HD2	1.92	0.52
1:A:727:ASN:CG	1:A:730:VAL:HG22	2.31	0.51
1:A:1003:ASP:OD2	1:A:1034:SER:OG	2.29	0.50
1:A:827:LYS:CG	1:A:1039:GLN:HG2	2.40	0.49
1:A:540:ARG:CZ	1:A:605:VAL:HG11	2.42	0.49
1:A:817:ASN:OD1	1:A:965:PRO:HD3	2.13	0.49
1:A:1054:GLN:CD	3:A:1346:HOH:O	2.52	0.48
1:A:793:TRP:HE3	1:A:903:MET:HG2	1.79	0.48
1:A:830:VAL:O	1:A:831:THR:HG22	2.14	0.48
1:A:834:ASN:HB2	1:A:1014:SER:OG	2.14	0.47
1:A:831:THR:CG2	1:A:1038:GLN:O	2.62	0.47
1:A:1019:ASN:HB3	1:A:1034:SER:HB2	1.97	0.47
1:A:984:VAL:HG21	1:A:991:ARG:CZ	2.46	0.45
1:A:924:LYS:HB2	1:A:930:TYR:CD2	2.51	0.45
1:A:942:ILE:H	1:A:942:ILE:HD13	1.81	0.45
1:A:903:MET:HA	1:A:903:MET:HE2	1.99	0.44
1:A:830:VAL:C	1:A:831:THR:CG2	2.85	0.44
1:A:899:ILE:HG13	1:A:900:LEU:HG	1.99	0.44
1:A:879:LYS:O	1:A:883:ASP:HA	2.17	0.43
1:A:1028:LYS:HE3	3:A:1310:HOH:O	2.18	0.43
1:A:564:ASP:HB3	1:A:569:LEU:HD11	2.01	0.43
1:A:976:TYR:CE2	1:A:978:GLY:HA2	2.54	0.42
1:A:539:PRO:O	1:A:543:ILE:HG13	2.19	0.42
1:A:903:MET:HB2	1:A:903:MET:HE3	1.82	0.42
1:A:1005:LYS:NZ	3:A:1337:HOH:O	2.53	0.42
1:A:821:ASP:OD2	1:A:821:ASP:N	2.50	0.42
1:A:885:ASP:HB2	3:A:1316:HOH:O	2.20	0.41
1:A:632:TRP:O	1:A:636:GLN:HG2	2.20	0.41
1:A:825:ILE:HG12	1:A:971:GLU:OE1	2.20	0.41
1:A:976:TYR:CZ	1:A:978:GLY:HA2	2.56	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:891:LEU:O	1:A:915:GLN:HA	2.21	0.41
1:A:924:LYS:O	1:A:925:SER:CB	2.69	0.41
1:A:846:LEU:HD22	1:A:1057:LEU:HD21	2.03	0.40
1:A:834:ASN:HB3	1:A:1036:ASN:HD22	1.81	0.40
1:A:871:LEU:HD23	1:A:871:LEU:HA	1.93	0.40
1:A:1054:GLN:HG3	3:A:1346:HOH:O	2.21	0.40
1:A:945:LYS:HE3	3:A:1472:HOH:O	2.20	0.40
1:A:829:TYR:OH	1:A:1041:LYS:HE2	2.21	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	539/573 (94%)	524 (97%)	12 (2%)	3 (1%)	25 10

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	924	LYS
1	A	926	ASP
1	A	613	ARG

### 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	490/521 (94%)	478 (98%)	12 (2%)	49 26

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

Mol	Chain	Res	Type
1	A	540	ARG
1	A	644	TRP
1	A	821	ASP
1	A	834	ASN
1	A	844	MET
1	A	903	MET
1	A	922	MET
1	A	923	GLN
1	A	942	ILE
1	A	1021	ASN
1	A	1054	GLN
1	A	1073	PRO

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

Mol	Chain	Res	Type
1	A	690	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 carbohydrates in this entry.

## 5.6 Ligand geometry [i](#)

1 ligand is 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	GOL	A	1200	-	5,5,5	0.33	0	5,5,5	0.49	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	GOL	A	1200	-	-	2/4/4/4	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	1200	GOL	C1-C2-C3-O3
2	A	1200	GOL	O2-C2-C3-O3

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	539/573 (94%)	0.40	48 (8%) <b>9</b> <b>12</b>	12, 23, 47, 69	0

All (48) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	538	ALA	8.4
1	A	612	LYS	7.9
1	A	833	ASN	6.2
1	A	823	VAL	6.1
1	A	832	MET	6.0
1	A	681	HIS	5.9
1	A	1076	PHE	5.6
1	A	923	GLN	5.6
1	A	540	ARG	4.9
1	A	682	HIS	4.8
1	A	1075	ASP	4.6
1	A	884	ILE	4.5
1	A	683	ASN	4.3
1	A	541	ASN	3.8
1	A	830	VAL	3.8
1	A	1000	LEU	3.8
1	A	821	ASP	3.7
1	A	684	GLY	3.7
1	A	685	LYS	3.6
1	A	614	ARG	3.5
1	A	539	PRO	3.5
1	A	707	ALA	3.4
1	A	709	ASP	3.4
1	A	922	MET	3.3
1	A	1074	SER	3.3
1	A	742	GLN	3.2
1	A	924	LYS	3.0

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Mol	Chain	Res	Type	RSRZ
1	A	759	ILE	3.0
1	A	913	THR	2.9
1	A	891	LEU	2.9
1	A	822	ASN	2.8
1	A	826	LYS	2.8
1	A	831	THR	2.7
1	A	893	GLY	2.7
1	A	901	GLU	2.7
1	A	963	VAL	2.7
1	A	834	ASN	2.6
1	A	824	GLN	2.6
1	A	926	ASP	2.5
1	A	866	ILE	2.4
1	A	609	LYS	2.3
1	A	902	GLN	2.2
1	A	574	ILE	2.2
1	A	911	TRP	2.2
1	A	974	PHE	2.1
1	A	761	TRP	2.1
1	A	885	ASP	2.1
1	A	987	GLU	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 carbohydrates 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	GOL	A	1200	6/6	0.94	0.10	21,28,39,39	0

## 6.5 Other polymers

There are no such residues in this entry.