Full wwPDB X-ray Structure Validation Report ⓘ

May 28, 2020 – 10:05 pm BST

PDB ID	:	2IL5
Title	:	Structure of Protein of Unknown Function SA2116 from Staphylococcus aureus
Authors	:	Osipiuk, J.; Quartey, P.; Holzle, D.; Joachimiak, A.; Midwest Center for Structural Genomics (MCSG)
Deposited on	:	2006-10-02
Resolution	:	2.30 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 following versions of software and data (see references ⓘ) were used in the production of this report:

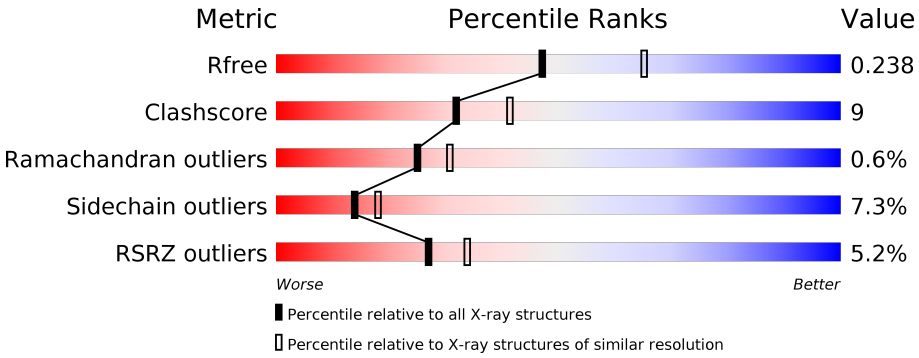
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.11
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.11

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.30 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	5042 (2.30-2.30)
Clashscore	141614	5643 (2.30-2.30)
Ramachandran outliers	138981	5575 (2.30-2.30)
Sidechain outliers	138945	5575 (2.30-2.30)
RSRZ outliers	127900	4938 (2.30-2.30)

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 >=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 <=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	171	5%, 74%, 17%, • • 5%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 1455 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 Hypothetical protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	Se			
1	A	162	1372	874	223	267	8	0	1	0

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-2	SER	-	CLONING ARTIFACT	UNP Q2FEH1
A	-1	ASN	-	CLONING ARTIFACT	UNP Q2FEH1
A	0	ALA	-	CLONING ARTIFACT	UNP Q2FEH1
A	1	MSE	MET	MODIFIED RESIDUE	UNP Q2FEH1
A	41	MSE	MET	MODIFIED RESIDUE	UNP Q2FEH1
A	83	MSE	MET	MODIFIED RESIDUE	UNP Q2FEH1
A	89	MSE	MET	MODIFIED RESIDUE	UNP Q2FEH1
A	93	MSE	MET	MODIFIED RESIDUE	UNP Q2FEH1
A	117	MSE	MET	MODIFIED RESIDUE	UNP Q2FEH1
A	147	MSE	MET	MODIFIED RESIDUE	UNP Q2FEH1
A	154	MSE	MET	MODIFIED RESIDUE	UNP Q2FEH1
A	158	MSE	MET	MODIFIED RESIDUE	UNP Q2FEH1

- Molecule 2 is SODIUM ION (three-letter code: NA) (formula: Na).

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

- Molecule 3 is water.

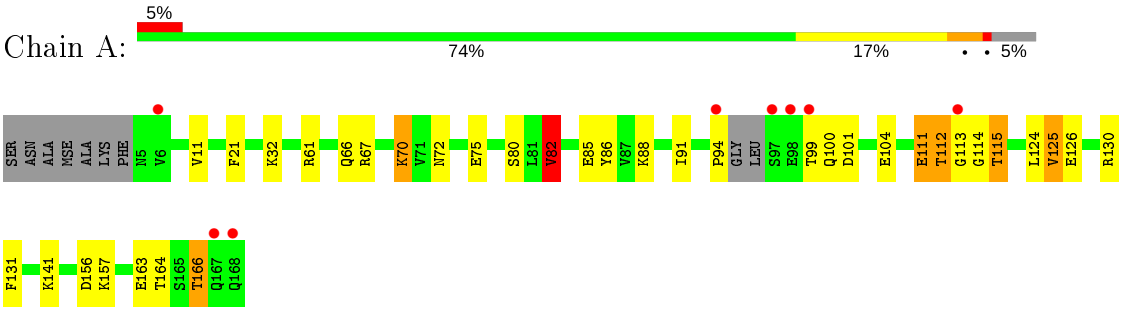
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
3	A	82	82	82	0	0

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: Hypothetical protein

Chain A:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 62	Depositor
Cell constants a, b, c, α, β, γ	107.59Å 107.59Å 39.10Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	35.20 – 2.30 35.22 – 2.30	Depositor EDS
% Data completeness (in resolution range)	98.4 (35.20-2.30) 98.4 (35.22-2.30)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	1.88 (at 2.29Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.202 , 0.245 0.193 , 0.238	Depositor DCC
R_{free} test set	1121 reflections (9.72%)	wwPDB-VP
Wilson B-factor (Å^2)	49.8	Xtriage
Anisotropy	0.503	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.32 , 46.6	EDS
L-test for twinning[2]	$< \|L\| > = 0.49$, $< L^2 > = 0.33$	Xtriage
Estimated twinning fraction	0.039 for h,-h-k,-l	Xtriage
F_o,F_c correlation	0.95	EDS
Total number of atoms	1455	wwPDB-VP
Average B, all atoms (Å^2)	49.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.11% of the height of the origin peak. No significant pseudotranslation is detected.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ 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: NA

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	0.81	0/1393	0.85	4/1853 (0.2%)

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.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	130	ARG	NE-CZ-NH2	-8.22	116.19	120.30
1	A	130	ARG	NE-CZ-NH1	6.58	123.59	120.30
1	A	156	ASP	CB-CA-C	5.45	121.29	110.40
1	A	82	VAL	CB-CA-C	5.13	121.15	111.40

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	111	GLU	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	1372	0	1335	24	0
2	A	1	0	0	0	0
3	A	82	0	0	4	0
All	All	1455	0	1335	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 9.

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:112:THR:HG23	1:A:113:GLY:H	1.33	0.92
1:A:21:PHE:O	1:A:114:GLY:HA3	1.74	0.88
1:A:114:GLY:HA2	3:A:267:HOH:O	1.85	0.75
1:A:91:ILE:HB	1:A:101:ASP:HB2	1.73	0.69
1:A:21:PHE:O	1:A:115:THR:HG22	2.02	0.60
1:A:112:THR:HG23	1:A:113:GLY:N	2.12	0.57
1:A:82:VAL:HG22	1:A:85:GLU:HB3	1.89	0.55
1:A:157:LYS:HD2	3:A:275:HOH:O	2.07	0.55
1:A:111:GLU:O	1:A:112:THR:HG22	2.08	0.53
1:A:101:ASP:OD1	3:A:205:HOH:O	2.18	0.53
1:A:82:VAL:HG13	1:A:86:TYR:HB3	1.92	0.52
1:A:94:PRO:HG3	1:A:100:GLN:NE2	2.25	0.52
1:A:61:ARG:HG3	1:A:75:GLU:HG2	1.93	0.50
1:A:67:ARG:O	1:A:70:LYS:HG3	2.12	0.49
1:A:112:THR:CG2	1:A:113:GLY:H	2.07	0.47
1:A:99:THR:HG21	1:A:124:LEU:HB2	1.97	0.46
1:A:66:GLN:HB3	3:A:218:HOH:O	2.14	0.46
1:A:67:ARG:O	1:A:70:LYS:HE3	2.16	0.45
1:A:11:VAL:HG13	1:A:125:VAL:CG1	2.48	0.44
1:A:112:THR:CG2	1:A:113:GLY:N	2.78	0.44
1:A:88:LYS:HG3	1:A:104:GLU:HG2	2.00	0.43
1:A:163:GLU:O	1:A:166:THR:HG22	2.19	0.42
1:A:80:SER:HB3	1:A:88:LYS:HB3	2.01	0.42
1:A:111:GLU:C	1:A:112:THR:HG22	2.40	0.41

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	159/171 (93%)	151 (95%)	7 (4%)	1 (1%)	25 31

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	112	THR

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	151/147 (103%)	140 (93%)	11 (7%)	14 18

All (11) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	32	LYS
1	A	70	LYS
1	A	72	ASN
1	A	82	VAL
1	A	115	THR
1	A	125	VAL
1	A	126	GLU
1	A	131	PHE
1	A	141	LYS
1	A	164	THR

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Mol	Chain	Res	Type
1	A	166	THR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	100	GLN
1	A	133	ASN
1	A	160	HIS

5.3.3 RNA

There are no RNA molecules in this entry.

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

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates

There are no carbohydrates in this entry.

5.6 Ligand geometry

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

There are no such residues in this entry.

5.8 Polymer linkage issues

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^{th} 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^2)	Q<0.9
1	A	154/171 (90%)	0.22	8 (5%) 27 34	38, 47, 67, 86	0

All (8) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	97	SER	6.2
1	A	167	GLN	4.1
1	A	168	GLN	4.0
1	A	98	GLU	3.2
1	A	113	GLY	3.0
1	A	6	VAL	2.8
1	A	99	THR	2.1
1	A	94	PRO	2.1

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

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates

There are no carbohydrates in this entry.

6.4 Ligands

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^{th} 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^2)	Q<0.9
2	NA	A	201	1/1	0.90	0.13	66,66,66,66	0

6.5 Other polymers (i)

There are no such residues in this entry.