



# Full wwPDB X-ray Structure Validation Report ⓘ

Mar 13, 2017 – 05:10 PM EDT

PDB ID : 5IHW  
Title : The crystal structure of SdrE from staphylococcus aureus  
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Deposited on : 2016-02-29  
Resolution : 1.25 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

<http://wwpdb.org/validation/2016/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  
Xtriage (Phenix) : 1.9-1692  
EDS : rb-20029077  
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)  
Refmac : 5.8.0135  
CCP4 : 6.5.0  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : rb-20029077

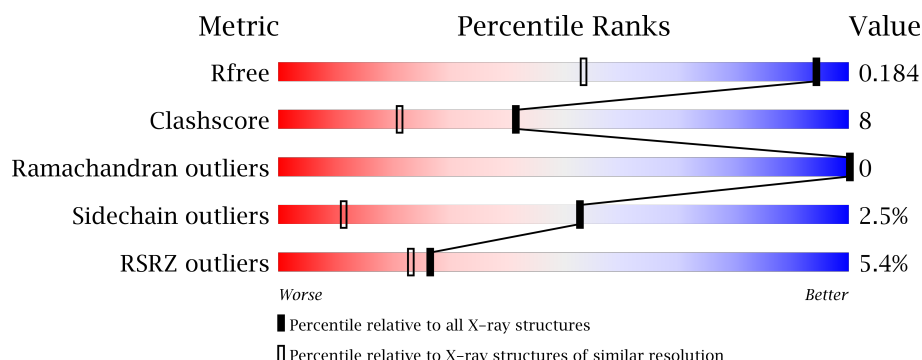
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 1.25 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))
$R_{free}$	100719	1690 (1.30-1.22)
Clashscore	112137	1781 (1.30-1.22)
Ramachandran outliers	110173	1716 (1.30-1.22)
Sidechain outliers	110143	1714 (1.30-1.22)
RSRZ outliers	101464	1693 (1.30-1.22)

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. 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	314	<div> <div>5%</div> <div>84%</div> <div>14%</div> <div>..</div> </div>

## 2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 2807 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 Serine-aspartate repeat-containing protein E.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	314	Total	C	N	O	S	0	1	0
			2488	1555	400	530	3			

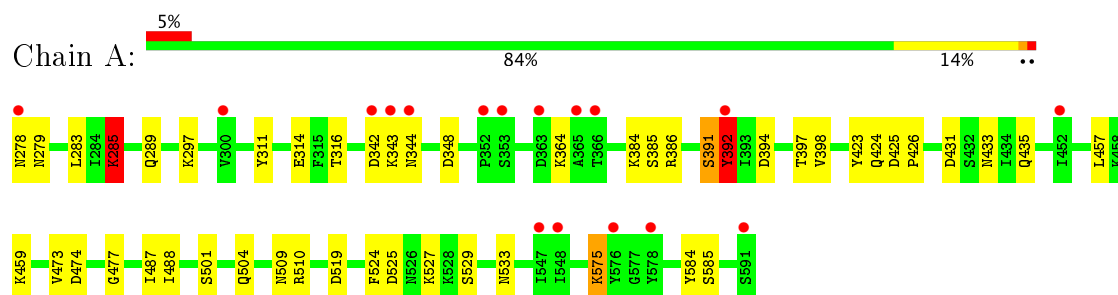
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	319	Total	O	0	0
			319	319		

### 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 ( $\text{RSRZ} > 2$ ). 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: Serine-aspartate repeat-containing protein E



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	39.26 Å   41.56 Å   51.14 Å 95.59°   97.55°   99.55°	Depositor
Resolution (Å)	50.32 – 1.25 22.67 – 1.25	Depositor EDS
% Data completeness (in resolution range)	87.0 (50.32-1.25) 83.3 (22.67-1.25)	Depositor EDS
$R_{merge}$	0.06	Depositor
$R_{sym}$	0.06	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	3.24 (at 1.25 Å)	Xtriage
Refinement program	REFMAC	Depositor
R, $R_{free}$	0.143   ,   0.176 0.153   ,   0.184	Depositor DCC
$R_{free}$ test set	2416 reflections (3.30%)	DCC
Wilson B-factor (Å <sup>2</sup> )	10.7	Xtriage
Anisotropy	0.263	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.44 , 60.3	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.97	EDS
Total number of atoms	2807	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	17.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 8.91% 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](#)

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	1.36	17/2528 (0.7%)	1.27	19/3429 (0.6%)

All (17) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	392[A]	TYR	CE1-CZ	-17.65	1.15	1.38
1	A	392[B]	TYR	CE1-CZ	-17.65	1.15	1.38
1	A	585	SER	CA-CB	9.93	1.67	1.52
1	A	385	SER	CB-OG	-9.61	1.29	1.42
1	A	392[A]	TYR	CB-CG	7.18	1.62	1.51
1	A	392[B]	TYR	CB-CG	7.18	1.62	1.51
1	A	504	GLN	CG-CD	6.36	1.65	1.51
1	A	391	SER	CB-OG	-6.14	1.34	1.42
1	A	392[A]	TYR	CD1-CE1	-5.86	1.30	1.39
1	A	392[B]	TYR	CD1-CE1	-5.86	1.30	1.39
1	A	392[A]	TYR	CG-CD1	-5.58	1.31	1.39
1	A	392[B]	TYR	CG-CD1	-5.58	1.31	1.39
1	A	529	SER	CB-OG	-5.24	1.35	1.42
1	A	423	TYR	CB-CG	5.06	1.59	1.51
1	A	585	SER	CB-OG	-5.03	1.35	1.42
1	A	426	PRO	N-CA	-5.02	1.38	1.47
1	A	423	TYR	CA-C	5.01	1.66	1.52

All (19) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	392[A]	TYR	CB-CG-CD2	16.07	130.64	121.00
1	A	392[B]	TYR	CB-CG-CD2	16.07	130.64	121.00
1	A	392[A]	TYR	OH-CZ-CE2	10.37	148.10	120.10
1	A	392[B]	TYR	OH-CZ-CE2	10.37	148.10	120.10
1	A	285	LYS	CD-CE-NZ	-10.35	87.89	111.70
1	A	392[A]	TYR	CE1-CZ-OH	-10.00	93.11	120.10
1	A	392[B]	TYR	CE1-CZ-OH	-10.00	93.11	120.10

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	392[A]	TYR	CG-CD1-CE1	9.28	128.72	121.30
1	A	392[B]	TYR	CG-CD1-CE1	9.28	128.72	121.30
1	A	348	ASP	CB-CG-OD2	-7.86	111.23	118.30
1	A	519	ASP	CB-CG-OD2	-7.59	111.47	118.30
1	A	392[A]	TYR	CB-CG-CD1	-7.50	116.50	121.00
1	A	392[B]	TYR	CB-CG-CD1	-7.50	116.50	121.00
1	A	425	ASP	CB-CG-OD1	-5.82	113.06	118.30
1	A	525	ASP	CB-CG-OD1	5.81	123.53	118.30
1	A	392[A]	TYR	CG-CD2-CE2	-5.21	117.13	121.30
1	A	392[B]	TYR	CG-CD2-CE2	-5.21	117.13	121.30
1	A	474	ASP	CB-CG-OD1	5.07	122.86	118.30
1	A	510	ARG	NE-CZ-NH1	-5.01	117.80	120.30

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	2488	0	2404	39	0
2	A	319	0	0	2	0
All	All	2807	0	2404	39	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 8.

All (39) 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:314:GLU:CD	1:A:386:ARG:NH1	2.14	1.00
1:A:314:GLU:CD	1:A:386:ARG:HH11	1.64	1.00
1:A:397:THR:HG21	2:A:693:HOH:O	1.66	0.96
1:A:314:GLU:OE2	1:A:386:ARG:NH1	2.01	0.94
1:A:394:ASP:CG	1:A:397:THR:HG23	1.90	0.92
1:A:278:ASN:HD22	1:A:279:ASN:H	1.26	0.83

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:575:LYS:H	1:A:575:LYS:HD2	1.45	0.81
1:A:392[A]:TYR:HD1	1:A:392[A]:TYR:N	1.77	0.80
1:A:342:ASP:OD1	1:A:343:LYS:HD3	1.85	0.77
1:A:278:ASN:ND2	1:A:279:ASN:H	1.85	0.74
1:A:392[A]:TYR:CD1	1:A:392[A]:TYR:N	2.51	0.74
1:A:391:SER:C	1:A:392[A]:TYR:CD1	2.63	0.71
1:A:524:PHE:HA	1:A:527:LYS:HG3	1.74	0.69
1:A:575:LYS:HD2	1:A:575:LYS:N	2.07	0.69
1:A:575:LYS:CD	1:A:575:LYS:H	2.08	0.67
1:A:394:ASP:OD2	1:A:397:THR:HG23	1.99	0.62
1:A:488:ILE:H	1:A:533:ASN:ND2	2.01	0.59
1:A:394:ASP:OD2	1:A:397:THR:CG2	2.51	0.58
1:A:391:SER:C	1:A:392[A]:TYR:HD1	2.02	0.57
1:A:433:ASN:HD21	1:A:457:LEU:H	1.54	0.56
1:A:314:GLU:HG3	2:A:777:HOH:O	2.06	0.55
1:A:424:GLN:HE21	1:A:509:ASN:HD22	1.55	0.54
1:A:316:THR:HG22	1:A:384:LYS:HG2	1.92	0.51
1:A:433:ASN:ND2	1:A:457:LEU:H	2.09	0.50
1:A:431:ASP:HA	1:A:459:LYS:HD2	1.93	0.49
1:A:314:GLU:CG	1:A:386:ARG:HH11	2.25	0.49
1:A:314:GLU:HG2	1:A:386:ARG:HD2	1.95	0.48
1:A:283:LEU:O	1:A:285:LYS:HE3	2.14	0.48
1:A:397:THR:OG1	1:A:398:VAL:HG23	2.14	0.47
1:A:488:ILE:H	1:A:533:ASN:HD21	1.63	0.46
1:A:424:GLN:HE21	1:A:509:ASN:ND2	2.15	0.45
1:A:364:LYS:HB2	1:A:364:LYS:HE3	1.78	0.43
1:A:285:LYS:HD2	1:A:316:THR:OG1	2.19	0.43
1:A:435:GLN:HE21	1:A:509:ASN:ND2	2.17	0.41
1:A:487:ILE:HA	1:A:533:ASN:HD21	1.85	0.41
1:A:314:GLU:CG	1:A:386:ARG:NH1	2.84	0.40
1:A:289:GLN:NE2	1:A:311:TYR:OH	2.55	0.40
1:A:394:ASP:CG	1:A:397:THR:CG2	2.77	0.40
1:A:473:VAL:CG1	1:A:477:GLY:HA2	2.52	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	313/314 (100%)	310 (99%)	3 (1%)	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	285/284 (100%)	277 (97%)	8 (3%)	49	9

All (8) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	285	LYS
1	A	297	LYS
1	A	344	ASN
1	A	392[A]	TYR
1	A	392[B]	TYR
1	A	501	SER
1	A	575	LYS
1	A	584	TYR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	278	ASN
1	A	289	GLN
1	A	334	ASN
1	A	405	ASN
1	A	433	ASN
1	A	509	ASN
1	A	533	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](#)

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	314/314 (100%)	0.21	17 (5%)	26 23	7, 15, 32, 54	0

All (17) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	343	LYS	4.9
1	A	300	VAL	4.5
1	A	576	TYR	4.0
1	A	365	ALA	3.6
1	A	392[A]	TYR	3.3
1	A	578	TYR	2.8
1	A	344	ASN	2.8
1	A	452	ILE	2.7
1	A	352	PRO	2.6
1	A	548	ILE	2.5
1	A	363	ASP	2.4
1	A	353	SER	2.3
1	A	342	ASP	2.2
1	A	591	SER	2.2
1	A	366	THR	2.1
1	A	278	ASN	2.0
1	A	547	ILE	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](#)

There are no ligands in this entry.

## 6.5 Other polymers [i](#)

There are no such residues in this entry.