



wwPDB X-ray Structure Validation Summary Report ⓘ

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PDB ID : 2NYD / pdb_00002nyd
Title : Crystal structure of Staphylococcus aureus hypothetical protein SA1388
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Deposited on : 2006-11-20
Resolution : 2.00 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

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>.

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

MolProbity : 4-5-2 with Phenix2.0
Xtrriage (Phenix) : 2.0
EDS : 3.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

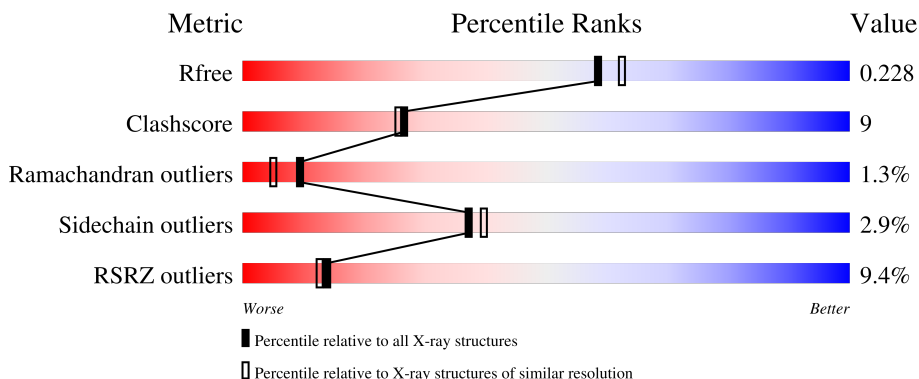
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.00 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}	180053	10052 (2.00-2.00)
Clashscore	190562	11152 (2.00-2.00)
Ramachandran outliers	187476	11031 (2.00-2.00)
Sidechain outliers	187428	11029 (2.00-2.00)
RSRZ outliers	180081	10067 (2.00-2.00)

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 ≥ 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	370	 5% 68% 11% • 19%
1	B	370	 12% 78% 15% •• 6%

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 5932 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 UPF0135 protein SA1388.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	300	2361	1507	388	454	12	0	1	0
1	B	348	2718	1739	441	524	14	0	0	0

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-3	ALA	-	cloning artifact	UNP P67273
A	-2	MET	-	cloning artifact	UNP P67273
A	-1	ASP	-	cloning artifact	UNP P67273
A	0	PRO	-	cloning artifact	UNP P67273
A	32	GLU	GLY	SEE REMARK 999	UNP P67273
A	115	VAL	ALA	SEE REMARK 999	UNP P67273
B	-3	ALA	-	cloning artifact	UNP P67273
B	-2	MET	-	cloning artifact	UNP P67273
B	-1	ASP	-	cloning artifact	UNP P67273
B	0	PRO	-	cloning artifact	UNP P67273
B	32	GLU	GLY	SEE REMARK 999	UNP P67273
B	115	VAL	ALA	SEE REMARK 999	UNP P67273

- Molecule 2 is ZINC ION (CCD ID: ZN) (formula: Zn).

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

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	419	Total 419	O 419	0	0
3	B	430	Total 430	O 430	0	0

4 Data and refinement statistics

Property	Value	Source
Space group	H 3	Depositor
Cell constants a, b, c, α , β , γ	132.55Å 132.55Å 125.67Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	84.51 – 2.00 84.51 – 2.00	Depositor EDS
% Data completeness (in resolution range)	95.0 (84.51-2.00) 95.0 (84.51-2.00)	Depositor EDS
R_{merge}	0.04	Depositor
R_{sym}	0.05	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.08 (at 2.00Å)	Xtrriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.173 , 0.231 0.172 , 0.228	Depositor DCC
R_{free} test set	2675 reflections (4.77%)	wwPDB-VP
Wilson B-factor (Å ²)	23.0	Xtrriage
Anisotropy	0.030	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 57.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.34$	Xtrriage
Estimated twinning fraction	0.017 for h,-h-k,-l	Xtrriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	5932	wwPDB-VP
Average B, all atoms (Å ²)	31.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.65% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical 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	0.98	1/2401 (0.0%)	0.97	3/3250 (0.1%)
1	B	0.97	2/2769 (0.1%)	0.95	2/3751 (0.1%)
All	All	0.98	3/5170 (0.1%)	0.96	5/7001 (0.1%)

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	B	0	1

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	276	ILE	CA-CB	6.60	1.62	1.54
1	A	255	ALA	CA-CB	6.56	1.63	1.53
1	B	34	VAL	CA-CB	5.10	1.60	1.54

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	288	GLY	N-CA-C	-6.18	106.96	113.58
1	A	73	THR	CB-CA-C	5.22	118.37	109.24
1	B	193	VAL	N-CA-C	5.10	114.64	106.88
1	A	222	THR	CA-C-N	5.03	126.13	119.84
1	A	222	THR	C-N-CA	5.03	126.13	119.84

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	219	PRO	Peptide

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	2361	0	2343	44	0
1	B	2718	0	2668	55	0
2	A	2	0	0	0	0
2	B	2	0	0	0	0
3	A	419	0	0	18	0
3	B	430	0	0	17	0
All	All	5932	0	5011	94	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.

The worst 5 of 94 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:346:ILE:HD13	3:A:678:HOH:O	1.10	1.23
1:B:346:ILE:CD1	3:B:832:HOH:O	1.87	1.18
1:B:346:ILE:HD13	3:B:832:HOH:O	1.46	1.06
1:B:33:ASP:HB3	3:B:698:HOH:O	1.59	1.03
1:B:55:GLU:HB3	3:B:689:HOH:O	1.58	1.02

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	295/370 (80%)	285 (97%)	9 (3%)	1 (0%)	36	35
1	B	344/370 (93%)	322 (94%)	15 (4%)	7 (2%)	6	2
All	All	639/740 (86%)	607 (95%)	24 (4%)	8 (1%)	9	5

5 of 8 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	140	LYS
1	B	219	PRO
1	B	141	ASP
1	B	155	LEU
1	B	192	ASP

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	259/322 (80%)	251 (97%)	8 (3%)	35	37
1	B	294/322 (91%)	286 (97%)	8 (3%)	39	42
All	All	553/644 (86%)	537 (97%)	16 (3%)	37	40

5 of 16 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	158	GLU
1	B	147	LYS
1	B	6	LEU
1	B	140	LYS
1	A	347	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 18 such sidechains are listed below:

Mol	Chain	Res	Type
1	B	206	GLN
1	B	320	ASN
1	B	218	HIS
1	B	24	ASN
1	B	160	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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 4 ligands modelled in this entry, 4 are 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 [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, 95th 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	300/370 (81%)	-0.11	18 (6%) 27 26	12, 21, 80, 80	1 (0%)
1	B	348/370 (94%)	0.11	43 (12%) 8 7	11, 21, 76, 80	0
All	All	648/740 (87%)	0.01	61 (9%) 14 13	11, 21, 78, 80	1 (0%)

The worst 5 of 61 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	220	TYR	6.9
1	A	131	TYR	5.5
1	B	-3	ALA	5.0
1	A	224	VAL	4.9
1	B	222	THR	4.8

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 oligosaccharides 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, 95th 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(\AA^2)	Q<0.9
2	ZN	A	401	1/1	0.97	0.14	50,50,50,50	0
2	ZN	B	402	1/1	0.97	0.17	47,47,47,47	0
2	ZN	B	401	1/1	0.98	0.14	41,41,41,41	0
2	ZN	A	402	1/1	0.98	0.14	48,48,48,48	0

6.5 Other polymers [i](#)

There are no such residues in this entry.