



wwPDB X-ray Structure Validation Summary Report

Feb 28, 2014 – 12:09 PM GMT

PDB ID : 4AQ1
Title : Structure of the SbsB S-layer protein of *Geobacillus stearothermophilus* PV72p2 in complex with nanobody KB6
Authors : Baranova, E.; Remaut, H.
Deposited on : 2012-04-12
Resolution : 2.42 Å(reported)

This is a wwPDB validation summary report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at <http://wwpdb.org/ValidationPDFNotes.html>

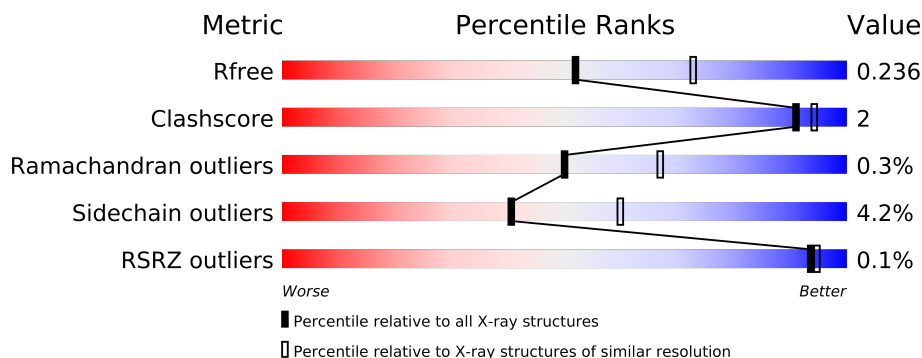
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity	:	4.02b-467
Mogul	:	1.15 2013
Xtriage (Phenix)	:	dev-1323
EDS	:	stable22639
Percentile statistics	:	21963
Refmac	:	5.8.0049
CCP4	:	6.3.0 (Settle)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP)	:	stable22683

1 Overall quality at a glance

The reported resolution of this entry is 2.42 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}	66092	2496 (2.44-2.40)
Clashscore	79885	3124 (2.44-2.40)
Ramachandran outliers	78287	3067 (2.44-2.40)
Sidechain outliers	78261	3068 (2.44-2.40)
RSRZ outliers	66119	2499 (2.44-2.40)

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. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	892	
2	B	130	
2	D	130	
3	C	892	

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 12335 atoms, of which 0 are hydrogen and 0 are deuterium.

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 SBSB PROTEIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	721	Total	C	N	O	S	82	0	0
			5363	3356	881	1124	2			

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	31	MET	-	EXPRESSION TAG	UNP Q45664
A	86	GLY	ASP	CONFLICT	UNP Q45664
A	139	ALA	ARG	CONFLICT	UNP Q45664
A	140	HIS	TYR	CONFLICT	UNP Q45664
A	500	GLU	ASP	CONFLICT	UNP Q45664
A	812	GLU	LYS	CONFLICT	UNP Q45664
A	921	GLY	-	EXPRESSION TAG	UNP Q45664
A	922	SER	-	EXPRESSION TAG	UNP Q45664

- Molecule 2 is a protein called NBKB6.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	124	Total	C	N	O	S	2	0	0
			924	572	163	184	5			
2	D	124	Total	C	N	O	S	7	0	0
			924	572	163	184	5			

- Molecule 3 is a protein called SBSB PROTEIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	627	Total	C	N	O	S	26	0	0
			4641	2899	761	979	2			

There are 9 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	31	MET	-	EXPRESSION TAG	UNP Q45664
C	86	GLY	ASP	CONFLICT	UNP Q45664
C	139	ALA	ARG	CONFLICT	UNP Q45664
C	140	HIS	TYR	CONFLICT	UNP Q45664
C	402	THR	SER	CONFLICT	UNP Q45664
C	500	GLU	ASP	CONFLICT	UNP Q45664
C	812	GLU	LYS	CONFLICT	UNP Q45664
C	921	GLY	-	EXPRESSION TAG	UNP Q45664
C	922	SER	-	EXPRESSION TAG	UNP Q45664

- Molecule 4 is CALCIUM ION (three-letter code: CA) (formula: Ca).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	4	Total Ca 4 4	0	0
4	C	4	Total Ca 4 4	0	0

- Molecule 5 is water.

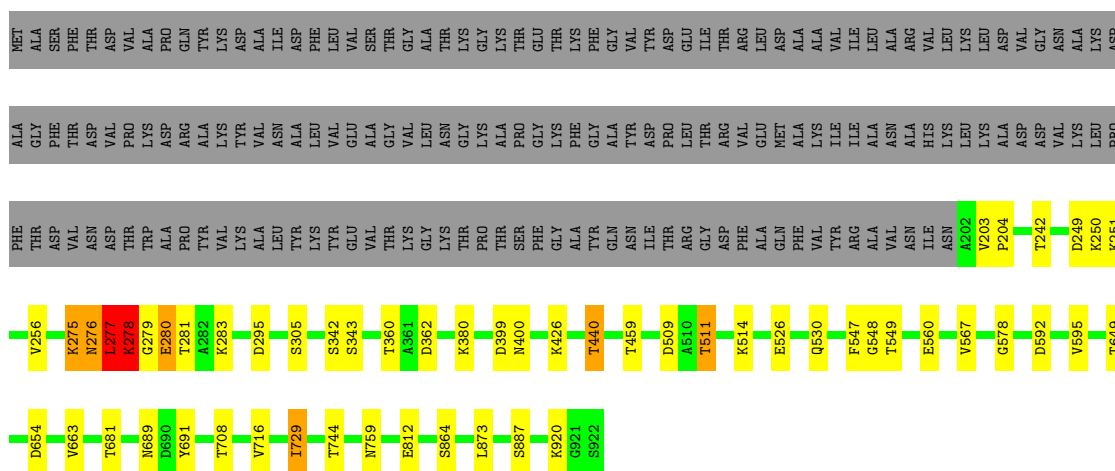
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	253	Total O 253 253	0	0
5	B	21	Total O 21 21	0	0
5	C	173	Total O 173 173	0	0
5	D	28	Total O 28 28	0	0

3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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: SBSB PROTEIN

Chain A:



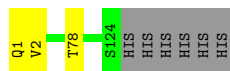
• Molecule 2: NBKB6

Chain B:



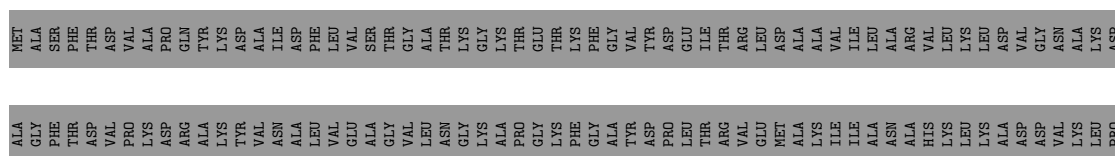
• Molecule 2: NBKB6

Chain D:



• Molecule 3: SBSB PROTEIN

Chain C:





4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	68.50Å 88.50Å 95.20Å 112.50° 101.60° 84.70°	Depositor
Resolution (Å)	43.81 – 2.42 43.81 – 2.42	Depositor EDS
% Data completeness (in resolution range)	97.0 (43.81-2.42) 97.1 (43.81-2.42)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.28 (at 2.42Å)	Xtriage
Refinement program	REFMAC 5.6.0117	Depositor
R, R_{free}	0.183 , 0.236 0.186 , 0.236	Depositor DCC
R_{free} test set	3729 reflections (5.27%)	DCC
Wilson B-factor (Å ²)	34.4	Xtriage
Anisotropy	0.056	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 20.2	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 74479 reflections	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	12335	wwPDB-VP
Average B, all atoms (Å ²)	39.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 11.66% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

5 Model quality ⓘ

5.1 Standard geometry ⓘ

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

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	0.67	3/5430 (0.1%)	0.77	1/7406 (0.0%)
2	B	0.70	1/943 (0.1%)	0.71	0/1271
2	D	0.70	0/943	0.76	0/1271
3	C	0.63	1/4698 (0.0%)	0.74	1/6409 (0.0%)
All	All	0.66	5/12014 (0.0%)	0.75	2/16357 (0.0%)

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

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	280	GLU	CA-CB	-10.73	1.30	1.53
1	A	275	LYS	CG-CD	-6.83	1.29	1.52
2	B	36	TRP	CD2-CE2	5.83	1.48	1.41
1	A	278	LYS	CD-CE	5.11	1.64	1.51
3	C	833	TRP	CD2-CE2	5.09	1.47	1.41

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	592	ASP	CB-CG-OD1	5.21	122.99	118.30
3	C	523	GLY	N-CA-C	-5.11	100.34	113.10

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	548	GLY	Peptide

5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	5363	0	65	16	0
2	B	924	0	0	3	0
2	D	924	0	10	1	0
3	C	4641	0	26	4	0
4	A	4	0	0	0	0
4	C	4	0	0	0	0
5	A	253	0	0	3	0
5	B	21	0	0	1	0
5	C	173	0	0	1	0
5	D	28	0	0	0	0
All	All	12335	0	101	23	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 2.

The worst 5 of 23 close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:440:THR:CG2	5:A:2087:HOH:O	2.28	0.81
1:A:511:THR:CG2	5:A:2024:HOH:O	2.38	0.71
1:A:276:ASN:O	1:A:277:LEU:C	2.39	0.60
1:A:295:ASP:N	1:A:295:ASP:OD1	2.36	0.58
1:A:547:PHE:O	1:A:578:GLY:O	2.22	0.56

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	719/892 (81%)	692 (96%)	23 (3%)	4 (1%)	33	46
2	B	122/130 (94%)	114 (93%)	8 (7%)	0	100	100
2	D	122/130 (94%)	119 (98%)	3 (2%)	0	100	100
3	C	625/892 (70%)	608 (97%)	16 (3%)	1 (0%)	56	74
All	All	1588/2044 (78%)	1533 (96%)	50 (3%)	5 (0%)	50	67

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	250	LYS
1	A	277	LEU
1	A	204	PRO
3	C	867	VAL
1	A	279	GLY

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	617/753 (82%)	586 (95%)	31 (5%)	34	51
2	B	92/98 (94%)	86 (94%)	6 (6%)	24	36
2	D	92/98 (94%)	91 (99%)	1 (1%)	84	94
3	C	534/753 (71%)	516 (97%)	18 (3%)	49	69
All	All	1335/1702 (78%)	1279 (96%)	56 (4%)	40	60

5 of 56 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	729	ILE
2	B	12	VAL
3	C	816	THR
1	A	744	THR
1	A	864	SER

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

5.3.3 RNA ⓘ

There are no RNA chains 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 8 ligands modelled in this entry, 8 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.

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, 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(Å²)	Q<0.9	
1	A	721/892 (80%)	-0.40	0	100	100	16, 34, 70, 109	26 (3%)
2	B	124/130 (95%)	-0.15	0	100	100	20, 45, 95, 127	1 (0%)
2	D	124/130 (95%)	-0.19	0	100	100	19, 36, 68, 86	3 (2%)
3	C	627/892 (70%)	-0.37	2 (0%)	91	92	18, 37, 65, 91	8 (1%)
All	All	1596/2044 (78%)	-0.35	2 (0%)	93	95	16, 36, 71, 127	38 (2%)

All (2) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
3	C	371	ASP	2.3
3	C	295	ASP	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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	RSR	LLDF	B-factors(\AA^2)	Q<0.9
4	CA	C	1922	1/1	0.08	-1.52	28,28,28,28	0
4	CA	C	1925	1/1	0.07	-1.80	33,33,33,33	0
4	CA	A	1926	1/1	0.07	-1.89	27,27,27,27	0
4	CA	A	1923	1/1	0.05	-2.08	25,25,25,25	0
4	CA	C	1923	1/1	0.05	-2.67	29,29,29,29	0
4	CA	C	1924	1/1	0.06	-3.19	24,24,24,24	0
4	CA	A	1925	1/1	0.04	-3.52	22,22,22,22	0
4	CA	A	1924	1/1	0.05	-4.12	23,23,23,23	0

6.5 Other polymers ⓘ

There are no such residues in this entry.