



# Full wwPDB X-ray Structure Validation Report ⓘ

Sep 27, 2017 – 12:08 AM EDT

PDB ID : 1VLI  
Title : Crystal structure of Spore coat polysaccharide biosynthesis protein spsE (BSU37870) from *Bacillus subtilis* at 2.38 Å resolution  
Authors : Joint Center for Structural Genomics (JCSG)  
Deposited on : unknown  
Resolution : 2.38 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-20030345
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-20030345

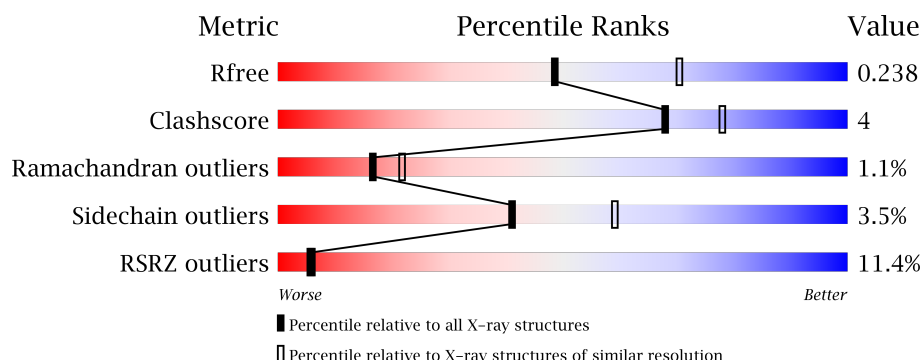
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.38 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	4388 (2.40-2.36)
Clashscore	112137	4984 (2.40-2.36)
Ramachandran outliers	110173	4907 (2.40-2.36)
Sidechain outliers	110143	4909 (2.40-2.36)
RSRZ outliers	101464	4423 (2.40-2.36)

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	385	<div> <div>10%</div> <div>81%</div> <div>11%</div> <div>7%</div> </div>

## 2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 2823 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 Spore coat polysaccharide biosynthesis protein spsE.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
			Total	C	N	O	S	Se			
1	A	358	2739	1752	457	518	4	8	0	0	0

There are 21 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-11	MSE	-	LEADER SEQUENCE	UNP P39625
A	-10	GLY	-	LEADER SEQUENCE	UNP P39625
A	-9	SER	-	LEADER SEQUENCE	UNP P39625
A	-8	ASP	-	LEADER SEQUENCE	UNP P39625
A	-7	LYS	-	LEADER SEQUENCE	UNP P39625
A	-6	ILE	-	LEADER SEQUENCE	UNP P39625
A	-5	HIS	-	LEADER SEQUENCE	UNP P39625
A	-4	HIS	-	LEADER SEQUENCE	UNP P39625
A	-3	HIS	-	LEADER SEQUENCE	UNP P39625
A	-2	HIS	-	LEADER SEQUENCE	UNP P39625
A	-1	HIS	-	LEADER SEQUENCE	UNP P39625
A	0	HIS	-	LEADER SEQUENCE	UNP P39625
A	1	MSE	MET	MODIFIED RESIDUE	UNP P39625
A	53	MSE	MET	MODIFIED RESIDUE	UNP P39625
A	59	MSE	MET	MODIFIED RESIDUE	UNP P39625
A	84	MSE	MET	MODIFIED RESIDUE	UNP P39625
A	86	MSE	MET	MODIFIED RESIDUE	UNP P39625
A	149	MSE	MET	MODIFIED RESIDUE	UNP P39625
A	179	MSE	MET	MODIFIED RESIDUE	UNP P39625
A	199	MSE	MET	MODIFIED RESIDUE	UNP P39625
A	260	MSE	MET	MODIFIED RESIDUE	UNP P39625

- Molecule 2 is ZINC ION (three-letter code: ZN) (formula: Zn).

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

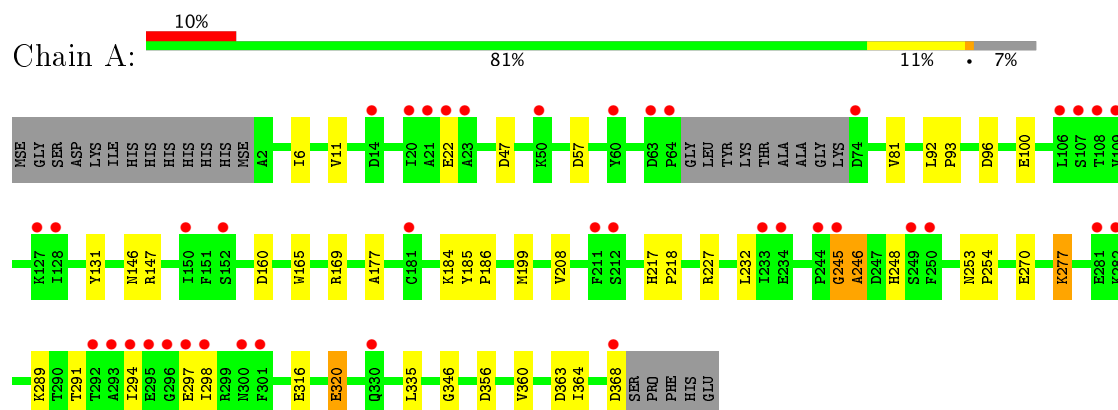
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	83	Total 83	O 83	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: Spore coat polysaccharide biosynthesis protein spsE



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 32 2 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	70.58 Å 70.58 Å 205.36 Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	29.29 – 2.38 29.29 – 2.38	Depositor EDS
% Data completeness (in resolution range)	99.8 (29.29-2.38) 99.9 (29.29-2.38)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	0.08	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.06 (at 2.39 Å)	Xtriage
Refinement program	REFMAC 5.2.0005	Depositor
R, $R_{free}$	0.190 , 0.238 0.189 , 0.238	Depositor DCC
$R_{free}$ test set	1256 reflections (5.38%)	DCC
Wilson B-factor (Å <sup>2</sup> )	49.4	Xtriage
Anisotropy	0.421	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.30 , 53.3	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.48$ , $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	0.046 for -h,-k,l	Xtriage
$F_o, F_c$ correlation	0.96	EDS
Total number of atoms	2823	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	63.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.00% 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

### 5.1 Standard geometry

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	0.90	4/2792 (0.1%)	0.83	5/3779 (0.1%)

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	270	GLU	CD-OE1	23.64	1.51	1.25
1	A	270	GLU	CD-OE2	15.27	1.42	1.25
1	A	270	GLU	CB-CG	6.60	1.64	1.52
1	A	316	GLU	CD-OE2	5.34	1.31	1.25

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	363	ASP	CB-CG-OD2	6.99	124.59	118.30
1	A	270	GLU	OE1-CD-OE2	6.04	130.54	123.30
1	A	47	ASP	CB-CG-OD2	5.89	123.60	118.30
1	A	160	ASP	CB-CG-OD2	5.85	123.57	118.30
1	A	368	ASP	CB-CG-OD2	5.11	122.90	118.30

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	2739	0	2690	24	0
2	A	1	0	0	0	0
3	A	83	0	0	4	0
All	All	2823	0	2690	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 4.

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:184:LYS:HG2	1:A:294:ILE:HD12	1.73	0.71
1:A:165:TRP:CZ2	1:A:169:ARG:HD2	2.26	0.71
1:A:291:THR:HG21	1:A:297:GLU:H	1.56	0.70
1:A:199:MSE:HE1	1:A:289:LYS:HB2	1.80	0.62
1:A:184:LYS:CB	1:A:294:ILE:HD12	2.30	0.62
1:A:199:MSE:CE	1:A:289:LYS:HB2	2.33	0.58
1:A:6:ILE:HD12	1:A:11:VAL:HG21	1.86	0.57
1:A:184:LYS:CG	1:A:294:ILE:HD12	2.35	0.55
1:A:165:TRP:CZ2	1:A:169:ARG:CD	2.90	0.54
1:A:246:ALA:H	1:A:248:HIS:HD2	1.57	0.53
1:A:320:GLU:HB2	3:A:412:HOH:O	2.10	0.50
1:A:227:ARG:HD2	1:A:277:LYS:HB3	1.94	0.49
1:A:92:LEU:HB2	1:A:93:PRO:HD3	1.98	0.46
1:A:245:GLY:HA2	1:A:248:HIS:HB2	1.98	0.45
1:A:346:GLY:HA2	3:A:441:HOH:O	2.17	0.44
1:A:217:HIS:HA	1:A:218:PRO:HD3	1.92	0.44
1:A:165:TRP:CE2	1:A:169:ARG:HD2	2.53	0.43
1:A:227:ARG:HD2	1:A:277:LYS:O	2.18	0.43
1:A:81:VAL:HB	3:A:438:HOH:O	2.18	0.43
1:A:131:TYR:HE1	1:A:298:ILE:HD12	1.84	0.43
1:A:185:TYR:HA	1:A:186:PRO:HA	1.85	0.42
1:A:253:ASN:HB2	1:A:254:PRO:CD	2.50	0.41
1:A:147:ARG:HG3	3:A:457:HOH:O	2.20	0.41
1:A:177:ALA:HA	1:A:208:VAL:O	2.21	0.41

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	354/385 (92%)	338 (96%)	12 (3%)	4 (1%)	17 22

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	246	ALA
1	A	320	GLU
1	A	277	LYS
1	A	245	GLY

### 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	283/305 (93%)	273 (96%)	10 (4%)	41 59

All (10) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	22	GLU
1	A	57	ASP
1	A	96	ASP
1	A	100	GLU
1	A	146	ASN
1	A	232	LEU
1	A	335	LEU

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Mol	Chain	Res	Type
1	A	356	ASP
1	A	360	VAL
1	A	364	ILE

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 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<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	350/385 (90%)	0.51	40 (11%) 6 6	48, 59, 89, 122	0

All (40) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	301	PHE	6.6
1	A	294	ILE	6.2
1	A	293	ALA	4.7
1	A	127	LYS	4.1
1	A	150	ILE	4.0
1	A	297	GLU	4.0
1	A	245	GLY	3.9
1	A	212	SER	3.9
1	A	296	GLY	3.8
1	A	128	ILE	3.7
1	A	295	GLU	3.7
1	A	181	CYS	3.6
1	A	20	ILE	3.6
1	A	108	THR	3.4
1	A	292	THR	3.2
1	A	368	ASP	3.2
1	A	63	ASP	3.1
1	A	234	GLU	3.1
1	A	60	TYR	2.9
1	A	281	GLU	2.9
1	A	106	LEU	2.9
1	A	107	SER	2.7
1	A	50	LYS	2.7
1	A	74	ASP	2.6
1	A	64	PRO	2.6
1	A	249	SER	2.6
1	A	300	ASN	2.5

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Mol	Chain	Res	Type	RSRZ
1	A	109	VAL	2.5
1	A	233	ILE	2.5
1	A	21	ALA	2.4
1	A	14	ASP	2.4
1	A	23	ALA	2.4
1	A	330	GLN	2.3
1	A	211	PHE	2.2
1	A	244	PRO	2.2
1	A	152	SER	2.2
1	A	250	PHE	2.1
1	A	22	GLU	2.1
1	A	282	LYS	2.1
1	A	298	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](#)

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, 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	LLDF	B-factors(Å <sup>2</sup> )	Q<0.9
2	ZN	A	374	1/1	0.92	0.31	-	57,57,57,57	1

## 6.5 Other polymers [i](#)

There are no such residues in this entry.