



# Full wwPDB X-ray Structure Validation Report ⓘ

Jul 2, 2025 – 12:08 AM JST

PDB ID : 9ILW / pdb\_00009ilw  
Title : Crystal structure of the periplasmic domain of sensor protein EvgS from Escherichia coli str. K-12 substr. MG1655  
Authors : Teng, Y.; Liu, R.; Luo, B.  
Deposited on : 2024-07-01  
Resolution : 2.70 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

<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 Phenix2.0rc1
Xtriage (Phenix)	:	2.0rc1
EDS	:	3.0
Percentile statistics	:	20231227.v01 (using entries in the PDB archive December 27th 2023)
CCP4	:	9.0.006 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.44

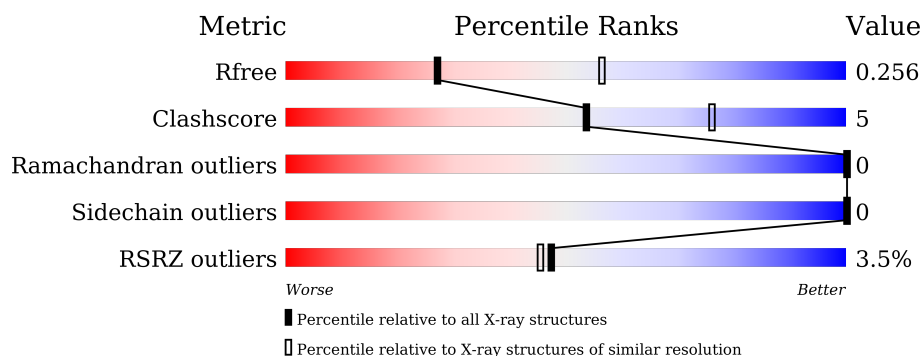
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.70 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))
$R_{free}$	164625	3333 (2.70-2.70)
Clashscore	180529	3684 (2.70-2.70)
Ramachandran outliers	177936	3633 (2.70-2.70)
Sidechain outliers	177891	3633 (2.70-2.70)
RSRZ outliers	164620	3333 (2.70-2.70)

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  $\geq 3$ , 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	513	 4% 86% 10% .
1	B	513	 3% 84% 12% .

## 2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 8024 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 Acid-sensing system histidine kinase EvgS.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	496	Total	C	N	O	S	0	0	0
			3964	2545	666	743	10			
1	B	496	Total	C	N	O	S	0	0	0
			3965	2548	665	742	10			

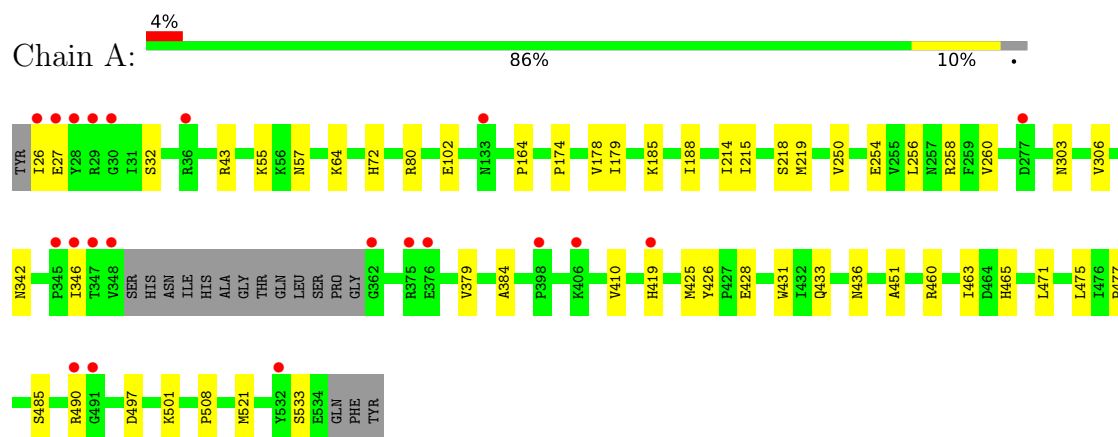
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	44	Total	O	0	0
			44	44		
2	B	51	Total	O	0	0
			51	51		

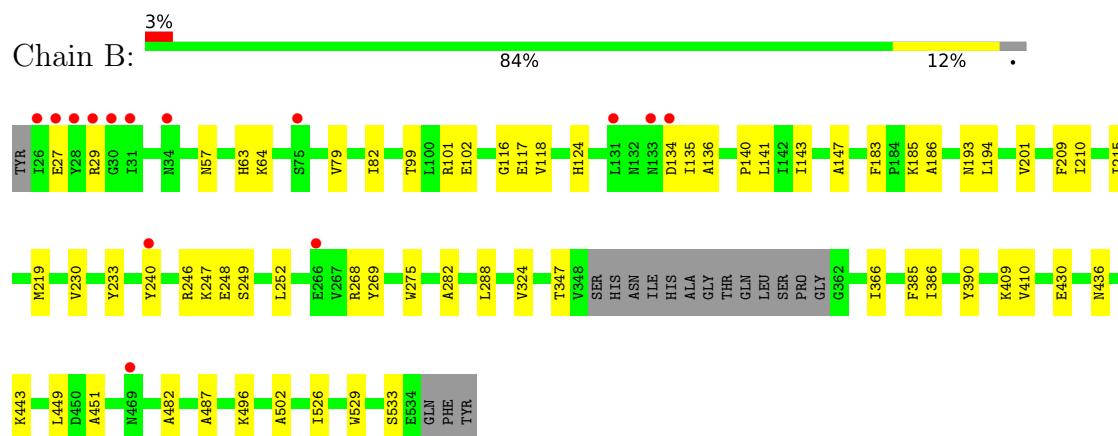
### 3 Residue-property plots

These plots are drawn for all protein, RNA, DNA and oligosaccharide 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: Acid-sensing system histidine kinase EvgS



- Molecule 1: Acid-sensing system histidine kinase EvgS



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	85.56Å 109.91Å 118.67Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	39.87 – 2.70 39.87 – 2.70	Depositor EDS
% Data completeness (in resolution range)	99.9 (39.87-2.70) 99.8 (39.87-2.70)	Depositor EDS
$R_{merge}$	0.25	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.57 (at 2.69Å)	Xtriage
Refinement program	PHENIX v1.0	Depositor
R, $R_{free}$	0.204 , 0.256 0.205 , 0.256	Depositor DCC
$R_{free}$ test set	1570 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	37.1	Xtriage
Anisotropy	0.483	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.35 , 51.5	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.93	EDS
Total number of atoms	8024	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	42.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.77% 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

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	0.12	0/4061	0.33	0/5540
1	B	0.14	0/4063	0.35	0/5543
All	All	0.13	0/8124	0.34	0/11083

There are no bond length outliers.

There are no bond angle outliers.

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	3964	0	3938	33	0
1	B	3965	0	3934	46	0
2	A	44	0	0	0	0
2	B	51	0	0	1	0
All	All	8024	0	7872	76	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 5.

All (76) 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:B:143:ILE:HG23	1:B:240:TYR:CE2	2.15	0.81

*Continued on next page...*

*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:64:LYS:HE2	1:A:72:HIS:HA	1.68	0.74
1:A:303:ASN:HD22	1:A:342:ASN:HB2	1.59	0.68
1:B:134:ASP:OD2	1:B:134:ASP:N	2.25	0.67
1:B:136:ALA:HB2	1:B:247:LYS:HG3	1.78	0.63
1:B:27:GLU:O	1:B:101:ARG:HA	2.00	0.61
1:B:143:ILE:HG23	1:B:240:TYR:HE2	1.65	0.60
1:B:193:ASN:ND2	2:B:601:HOH:O	2.32	0.60
1:A:174:PRO:HB2	1:A:178:VAL:HG11	1.85	0.58
1:A:303:ASN:ND2	1:A:342:ASN:HB2	2.18	0.58
1:B:246:ARG:HG3	1:B:248:GLU:HB2	1.83	0.58
1:B:324:VAL:HG23	1:B:347:THR:HG21	1.86	0.57
1:A:27:GLU:HB2	1:A:102:GLU:H	1.69	0.57
1:A:460:ARG:NH1	1:B:533:SER:O	2.37	0.56
1:B:410:VAL:HG22	1:B:451:ALA:HB3	1.87	0.55
1:A:410:VAL:HG22	1:A:451:ALA:HB3	1.89	0.55
1:A:250:VAL:O	1:A:254:GLU:HG2	2.07	0.54
1:B:249:SER:HB3	1:B:252:LEU:HB3	1.90	0.54
1:A:26:ILE:O	1:A:27:GLU:HG2	2.09	0.53
1:B:116:GLY:HA2	1:B:246:ARG:NH2	2.22	0.53
1:A:214:ILE:HG23	1:A:215:ILE:HG23	1.93	0.51
1:B:29:ARG:HH12	1:B:102:GLU:HB2	1.76	0.51
1:B:141:LEU:O	1:B:268:ARG:NH1	2.35	0.51
1:B:443:LYS:HB3	1:B:449:LEU:HD12	1.93	0.50
1:B:324:VAL:HG23	1:B:347:THR:CG2	2.42	0.49
1:A:43:ARG:O	1:A:258:ARG:NH1	2.47	0.48
1:A:379:VAL:O	1:A:490:ARG:HD2	2.13	0.48
1:A:306:VAL:O	1:A:346:ILE:HG22	2.14	0.47
1:B:230:VAL:HG11	1:B:233:TYR:CE1	2.49	0.47
1:B:101:ARG:NH2	1:B:117:GLU:HG2	2.29	0.47
1:B:134:ASP:OD2	1:B:135:ILE:N	2.49	0.46
1:B:82:ILE:HG12	1:B:275:TRP:CD1	2.51	0.45
1:B:269:TYR:HA	1:B:282:ALA:HB3	1.98	0.45
1:A:419:HIS:CD2	1:A:433:GLN:HG3	2.52	0.45
1:B:29:ARG:NH2	1:B:64:LYS:HD2	2.30	0.45
1:B:63:HIS:NE2	1:B:124:HIS:HB3	2.32	0.45
1:B:436:ASN:ND2	1:B:529:TRP:O	2.50	0.45
1:A:55:LYS:O	1:A:57:ASN:N	2.41	0.45
1:B:134:ASP:CG	1:B:135:ILE:HG12	2.41	0.45
1:B:288:LEU:HD11	1:B:502:ALA:HA	1.99	0.45
1:B:140:PRO:HB3	1:B:240:TYR:CD2	2.51	0.45
1:B:409:LYS:HG2	1:B:430:GLU:OE1	2.16	0.44

*Continued on next page...*

*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:385:PHE:CD2	1:B:386:ILE:HG13	2.53	0.44
1:B:496:LYS:HE2	1:B:496:LYS:HB3	1.91	0.44
1:A:179:ILE:HD12	1:A:188:ILE:HD13	1.99	0.44
1:A:463:ILE:HG12	1:A:471:LEU:HB2	2.00	0.44
1:A:508:PRO:HB3	1:B:526:ILE:HG23	2.00	0.43
1:B:194:LEU:HD23	1:B:219:MET:SD	2.59	0.43
1:B:390:TYR:CZ	1:B:482:ALA:HB2	2.53	0.43
1:A:497:ASP:O	1:A:501:LYS:HG3	2.18	0.43
1:A:164:PRO:HA	1:A:185:LYS:O	2.19	0.43
1:A:426:TYR:HB2	1:A:431:TRP:HZ2	1.82	0.43
1:A:256:LEU:O	1:A:260:VAL:HG23	2.19	0.43
1:B:185:LYS:HD3	1:B:185:LYS:HA	1.70	0.43
1:A:384:ALA:HA	1:A:485:SER:HB2	2.00	0.43
1:A:475:LEU:HD11	1:B:529:TRP:CH2	2.53	0.42
1:A:428:GLU:CD	1:A:428:GLU:H	2.24	0.42
1:A:436:ASN:HD22	1:A:533:SER:HB3	1.84	0.42
1:A:32:SER:HB3	1:A:80:ARG:CG	2.49	0.42
1:B:134:ASP:CG	1:B:135:ILE:N	2.78	0.42
1:B:366:ILE:HB	1:B:487:ALA:HB3	2.01	0.42
1:A:465:HIS:HE1	1:A:521:MET:O	2.03	0.42
1:B:29:ARG:HH22	1:B:64:LYS:HD2	1.85	0.42
1:B:118:VAL:O	1:B:246:ARG:NH2	2.53	0.42
1:B:147:ALA:O	1:B:210:ILE:HA	2.20	0.41
1:B:57:ASN:OD1	1:B:99:THR:HG23	2.20	0.41
1:A:32:SER:HB3	1:A:80:ARG:HG3	2.02	0.41
1:B:29:ARG:NH1	1:B:102:GLU:HB2	2.35	0.41
1:B:183:PHE:HB3	1:B:186:ALA:HB2	2.03	0.41
1:A:218:SER:OG	1:A:219:MET:N	2.54	0.41
1:B:29:ARG:HD3	1:B:79:VAL:HG23	2.02	0.41
1:A:379:VAL:C	1:A:490:ARG:CG	2.94	0.40
1:B:194:LEU:HD22	1:B:215:ILE:HG22	2.02	0.40
1:B:201:VAL:HG21	1:B:209:PHE:HB2	2.03	0.40
1:A:425:MET:HE1	1:A:477:PRO:O	2.20	0.40
1:A:32:SER:O	1:A:80:ARG:HG2	2.22	0.40

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	492/513 (96%)	476 (97%)	16 (3%)	0	100	100
1	B	492/513 (96%)	476 (97%)	16 (3%)	0	100	100
All	All	984/1026 (96%)	952 (97%)	32 (3%)	0	100	100

There are no Ramachandran outliers to report.

### 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	441/460 (96%)	441 (100%)	0	100	100
1	B	440/460 (96%)	440 (100%)	0	100	100
All	All	881/920 (96%)	881 (100%)	0	100	100

There are no protein residues with a non-rotameric sidechain to report.

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (15) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	153	HIS
1	A	166	ASN
1	A	213	ASN
1	A	253	ASN
1	A	285	ASN
1	A	296	GLN

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Res	Type
1	A	300	GLN
1	A	303	ASN
1	A	419	HIS
1	A	433	GLN
1	B	46	ASN
1	B	106	HIS
1	B	235	ASN
1	B	253	ASN
1	B	433	GLN

### 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](#)

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

### 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	496/513 (96%)	0.07	21 (4%) 41 39	21, 39, 74, 115	0
1	B	496/513 (96%)	0.03	14 (2%) 55 53	19, 37, 74, 102	0
All	All	992/1026 (96%)	0.05	35 (3%) 47 45	19, 38, 74, 115	0

All (35) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	28	TYR	5.5
1	A	26	ILE	5.4
1	A	28	TYR	4.9
1	B	26	ILE	4.6
1	B	30	GLY	4.6
1	A	347	THR	4.5
1	B	134	ASP	4.5
1	A	30	GLY	3.9
1	A	27	GLU	3.9
1	A	29	ARG	3.6
1	A	348	VAL	3.4
1	A	346	ILE	3.3
1	A	532	TYR	3.3
1	A	490	ARG	3.2
1	B	34	ASN	3.1
1	A	419	HIS	3.1
1	A	376	GLU	3.0
1	A	36	ARG	3.0
1	B	27	GLU	2.9
1	B	29	ARG	2.9
1	A	398	PRO	2.9
1	A	375	ARG	2.8
1	A	362	GLY	2.6
1	B	133	ASN	2.5

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Res	Type	RSRZ
1	B	131	LEU	2.4
1	B	75	SER	2.4
1	A	491	GLY	2.3
1	B	469	ASN	2.2
1	B	266	GLU	2.2
1	A	133	ASN	2.1
1	B	240	TYR	2.1
1	A	345	PRO	2.1
1	B	31	ILE	2.1
1	A	277	ASP	2.1
1	A	406	LYS	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 oligosaccharides 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.