



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 13, 2017 – 02:49 am GMT

PDB ID : 3I5B
Title : Crystal structure of the isolated GGDEF domain of WpsR from *Pseudomonas aeruginosa*
Authors : Navarro, M.V.A.S.; De, N.; Sondermann, H.
Deposited on : 2009-07-03
Resolution : 2.04 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

A user guide is available at

<http://wwpdb.org/validation/2016/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467
Mogul : 1.7.2 (RC1), CSD as538be (2017)
Xtriage (Phenix) : 1.9-1692
EDS : trunk28620
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) : recalc28949

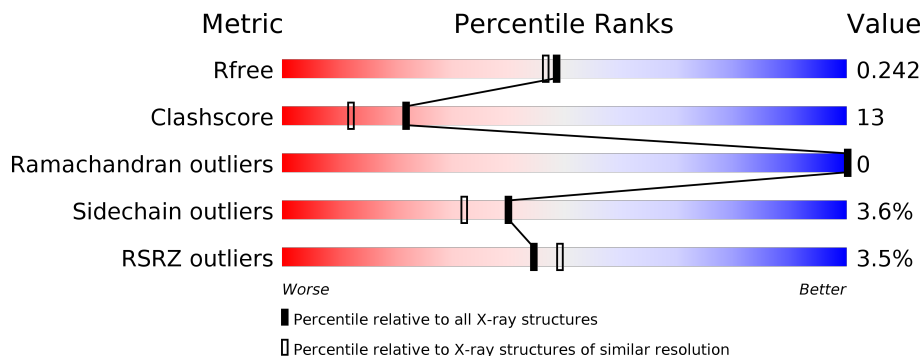
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.04 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	1316 (2.04-2.04)
Clashscore	112137	1394 (2.04-2.04)
Ramachandran outliers	110173	1383 (2.04-2.04)
Sidechain outliers	110143	1383 (2.04-2.04)
RSRZ outliers	101464	1319 (2.04-2.04)

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. 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	179	<div> <div>4%</div> <div> <div></div> <div>72%</div> <div>20%</div> <div>• 5%</div> </div> </div>
1	B	179	<div> <div>2%</div> <div> <div></div> <div>74%</div> <div>19%</div> <div>• 6%</div> </div> </div>

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 2855 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.

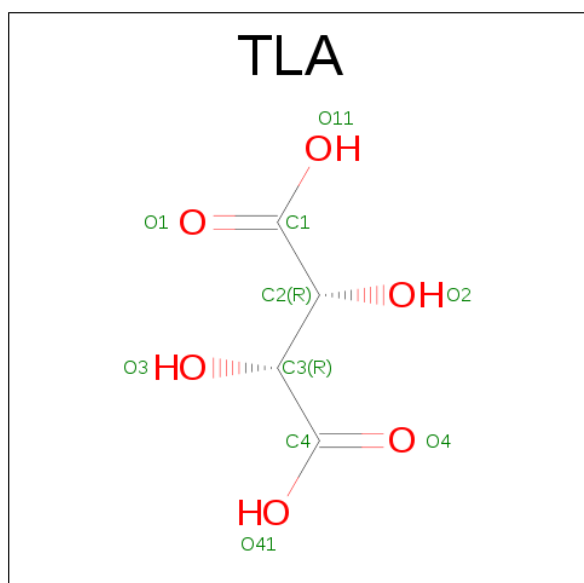
- Molecule 1 is a protein called WspR response regulator.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	170	Total	C	N	O	S	0	0	0
			1304	803	242	252	7			
1	B	169	Total	C	N	O	S	0	0	0
			1299	800	241	251	7			

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	169	ALA	-	EXPRESSION TAG	UNP Q9HXT9
A	170	LEU	-	EXPRESSION TAG	UNP Q9HXT9
A	171	MET	-	EXPRESSION TAG	UNP Q9HXT9
B	169	ALA	-	EXPRESSION TAG	UNP Q9HXT9
B	170	LEU	-	EXPRESSION TAG	UNP Q9HXT9
B	171	MET	-	EXPRESSION TAG	UNP Q9HXT9

- Molecule 2 is L(+)-TARTARIC ACID (three-letter code: TLA) (formula: C₄H₆O₆).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			10	4	6		
2	B	1	Total	C	O	0	0
			10	4	6		

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	106	Total	O	0	0
			106	106		
3	B	126	Total	O	0	0
			126	126		

4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	35.62Å 94.50Å 97.59Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	43.36 – 2.04 43.36 – 2.04	Depositor EDS
% Data completeness (in resolution range)	92.9 (43.36-2.04) 93.0 (43.36-2.04)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.11	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.69 (at 2.05Å)	Xtriage
Refinement program	PHENIX (phenix.refine)	Depositor
R, R_{free}	0.193 , 0.244 0.192 , 0.242	Depositor DCC
R_{free} test set	2034 reflections (9.97%)	DCC
Wilson B-factor (Å ²)	18.6	Xtriage
Anisotropy	0.262	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.40 , 53.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.55$, $\langle L^2 \rangle = 0.39$	Xtriage
Estimated twinning fraction	0.000 for -h,l,k	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	2855	wwPDB-VP
Average B, all atoms (Å ²)	20.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 54.13 % of the origin peak, indicating pseudo translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo translational symmetry is equal to 3.8151e-05. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

¹ Intensities estimated from amplitudes.

² 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: TLA

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	0.38	0/1324	0.52	0/1783
1	B	0.38	0/1319	0.52	0/1776
All	All	0.38	0/2643	0.52	0/3559

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	1304	0	1276	39	0
1	B	1299	0	1274	36	0
2	A	10	0	6	0	0
2	B	10	0	6	0	0
3	A	106	0	0	4	0
3	B	126	0	0	4	0
All	All	2855	0	2562	69	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 13.

All (69) 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:194:ARG:HD3	1:A:198:ARG:HH21	1.18	1.05
1:B:185:PHE:CE1	1:B:258:VAL:HG23	2.12	0.85
1:A:185:PHE:CE1	1:A:258:VAL:HG23	2.14	0.83
1:A:241:SER:HB2	3:A:352:HOH:O	1.80	0.81
1:A:237:ARG:HE	1:B:182:ARG:HH21	1.37	0.71
1:B:303:LEU:HD11	1:B:318:MET:SD	2.30	0.70
1:A:204:SER:OG	1:A:302:THR:HB	1.93	0.68
1:B:185:PHE:CZ	1:B:258:VAL:HG23	2.30	0.67
1:A:180:SER:OG	1:A:184:HIS:HD2	1.79	0.66
1:A:237:ARG:HE	1:B:182:ARG:NH2	1.93	0.66
1:A:210:VAL:CG1	1:A:294:LEU:HD21	2.26	0.65
1:B:264:PRO:HA	1:B:302:THR:HG21	1.79	0.64
1:A:234:GLY:O	1:A:238:GLU:HG3	1.98	0.64
1:A:264:PRO:HA	1:A:302:THR:HG21	1.82	0.61
1:A:227:GLU:OE1	1:A:230:ARG:NH1	2.34	0.61
1:B:246:LEU:HD12	1:B:247:ALA:N	2.15	0.61
1:B:169:ALA:N	3:B:156:HOH:O	2.33	0.60
1:B:204:SER:OG	1:B:302:THR:HB	2.03	0.59
1:A:337:MET:HG2	1:A:338:GLU:N	2.16	0.59
1:A:242:ARG:HG2	1:A:245:ASP:OD2	2.04	0.58
1:B:180:SER:OG	1:B:184:HIS:HD2	1.87	0.58
1:B:318:MET:HG2	1:B:337:MET:HE2	1.84	0.58
1:B:185:PHE:CZ	1:B:258:VAL:CG2	2.86	0.58
1:B:246:LEU:HD12	1:B:247:ALA:H	1.68	0.56
1:A:276:ARG:NH1	1:A:279:GLU:OE2	2.39	0.54
1:B:297:SER:HB2	1:B:327:LYS:HE2	1.88	0.54
1:B:182:ARG:NH1	1:B:250:TYR:CE1	2.75	0.54
1:B:242:ARG:HB2	1:B:245:ASP:OD2	2.08	0.54
1:A:325:GLN:HE22	1:A:336:LEU:H	1.55	0.53
1:A:325:GLN:NE2	1:A:336:LEU:H	2.07	0.53
1:A:170:LEU:HD23	1:B:321:GLN:HG3	1.91	0.52
1:A:268:ARG:HD3	1:A:336:LEU:HD11	1.91	0.52
1:B:325:GLN:HE22	1:B:336:LEU:H	1.60	0.49
1:A:337:MET:CG	1:A:338:GLU:N	2.75	0.49
1:B:182:ARG:HD3	1:B:250:TYR:CZ	2.48	0.49
1:B:259:LEU:HD11	1:B:270:LEU:HD22	1.94	0.49
1:A:294:LEU:HD23	3:A:42:HOH:O	2.12	0.48
1:B:236:ILE:HG23	1:B:257:MET:HE2	1.95	0.48
1:B:337:MET:HA	3:B:157:HOH:O	2.14	0.48
1:A:237:ARG:NE	1:B:182:ARG:HH21	2.09	0.47
1:A:242:ARG:HG2	1:A:245:ASP:CG	2.35	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:268:ARG:HD2	3:A:139:HOH:O	2.14	0.47
1:A:297:SER:HB2	1:A:327:LYS:HE2	1.97	0.47
1:A:231:GLN:HB3	1:A:281:LEU:HD13	1.97	0.46
1:A:337:MET:O	1:A:338:GLU:C	2.54	0.46
1:B:213:PHE:CE2	1:B:217:ASN:ND2	2.83	0.46
1:B:337:MET:HA	3:B:366:HOH:O	2.15	0.46
1:A:185:PHE:CZ	1:A:258:VAL:HG23	2.49	0.46
1:A:185:PHE:CZ	1:A:258:VAL:CG2	2.99	0.46
1:A:210:VAL:HG13	1:A:294:LEU:HD21	1.98	0.46
1:A:283:ILE:HB	1:A:294:LEU:HB2	1.98	0.45
1:A:170:LEU:HD21	1:B:317:GLU:HG3	1.99	0.45
1:B:325:GLN:NE2	1:B:336:LEU:H	2.15	0.45
1:A:169:ALA:N	3:A:361:HOH:O	2.49	0.45
1:A:170:LEU:HB3	1:B:320:ASP:OD2	2.17	0.44
1:A:194:ARG:CD	1:A:198:ARG:HH21	2.07	0.44
1:B:236:ILE:HG23	1:B:257:MET:CE	2.48	0.44
1:B:182:ARG:CD	1:B:250:TYR:CZ	3.01	0.43
1:B:240:CYS:HB2	1:B:245:ASP:HB2	2.00	0.43
1:A:217:ASN:ND2	1:A:222:HIS:HA	2.34	0.43
1:A:237:ARG:HA	1:A:240:CYS:HB2	2.01	0.42
1:A:300:VAL:HB	1:A:336:LEU:HD23	2.02	0.42
1:B:240:CYS:HA	3:B:121:HOH:O	2.19	0.42
1:A:210:VAL:HG13	1:A:294:LEU:CD2	2.49	0.42
1:A:213:PHE:CD2	1:A:253:GLU:HB3	2.56	0.41
1:B:234:GLY:O	1:B:238:GLU:HG3	2.21	0.41
1:B:276:ARG:NH1	1:B:279:GLU:OE2	2.52	0.41
1:B:240:CYS:HB2	1:B:245:ASP:CB	2.52	0.40
1:B:264:PRO:CA	1:B:302:THR:HG21	2.49	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	168/179 (94%)	164 (98%)	4 (2%)	0	100	100
1	B	167/179 (93%)	164 (98%)	3 (2%)	0	100	100
All	All	335/358 (94%)	328 (98%)	7 (2%)	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	137/144 (95%)	130 (95%)	7 (5%)	28	18
1	B	137/144 (95%)	134 (98%)	3 (2%)	57	51
All	All	274/288 (95%)	264 (96%)	10 (4%)	40	32

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

Mol	Chain	Res	Type
1	A	215	SER
1	A	240	CYS
1	A	242	ARG
1	A	289	ARG
1	A	294	LEU
1	A	302	THR
1	A	337	MET
1	B	289	ARG
1	B	302	THR
1	B	337	MET

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (14) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	184	HIS
1	A	217	ASN
1	A	287	GLN

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Mol	Chain	Res	Type
1	A	293	HIS
1	A	325	GLN
1	A	328	ASN
1	A	329	ASN
1	A	333	GLN
1	B	184	HIS
1	B	287	GLN
1	B	325	GLN
1	B	328	ASN
1	B	329	ASN
1	B	333	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 carbohydrates in this entry.

5.6 Ligand geometry [i](#)

2 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the chemical component dictionary. The Link column lists molecule types, if any, to which the group is linked. 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 $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
2	TLA	A	2	-	3,9,9	1.14	0	6,12,12	0.55	0
2	TLA	B	1	-	3,9,9	1.09	0	6,12,12	0.65	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the chemical component dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	TLA	A	2	-	-	0/4/12/12	0/0/0/0
2	TLA	B	1	-	-	0/4/12/12	0/0/0/0

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 [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	170/179 (94%)	0.21	8 (4%) 32 35	11, 18, 38, 60	0
1	B	169/179 (94%)	-0.00	4 (2%) 59 64	10, 18, 32, 56	0
All	All	339/358 (94%)	0.10	12 (3%) 44 49	10, 18, 36, 60	0

All (12) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	309	GLY	6.7
1	B	309	GLY	4.7
1	A	337	MET	4.1
1	A	308	GLY	4.1
1	B	337	MET	3.1
1	A	338	GLU	3.0
1	B	306	GLY	3.0
1	A	307	GLY	3.0
1	A	200	GLN	2.5
1	A	304	VAL	2.4
1	A	241	SER	2.2
1	B	307	GLY	2.2

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, 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	LLDF	B-factors(\AA^2)	Q<0.9
2	TLA	B	1	10/10	0.96	0.10	-0.49	10,14,16,19	0
2	TLA	A	2	10/10	0.97	0.09	-1.34	13,15,18,19	0

6.5 Other polymers [i](#)

There are no such residues in this entry.