



Full wwPDB X-ray Structure Validation Report ⓘ

May 18, 2020 – 07:39 am BST

PDB ID : 3KDG
Title : C-terminal domain of Bacillus subtilis MutL crystal form II
Authors : Guarne, A.; Pillon, M.C.
Deposited on : 2009-10-22
Resolution : 2.00 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

<https://www.wwpdb.org/validation/2017/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
Xtriage (Phenix)	:	1.13
EDS	:	2.11
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.11

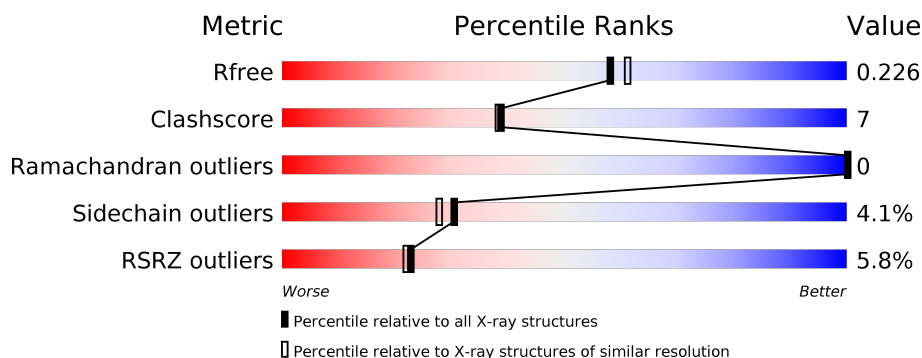
1 Overall quality at a glance

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}	130704	8085 (2.00-2.00)
Clashscore	141614	9178 (2.00-2.00)
Ramachandran outliers	138981	9054 (2.00-2.00)
Sidechain outliers	138945	9053 (2.00-2.00)
RSRZ outliers	127900	7900 (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 on 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	197	<div> <div>9%</div> <div> <div></div> <div>75%</div> <div>20%</div> <div>• •</div> </div> </div>
1	B	197	<div> <div>2%</div> <div> <div></div> <div>84%</div> <div>10%</div> <div>• 5%</div> </div> </div>

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 3294 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 DNA mismatch repair protein mutL.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	191	Total	C	N	O	S	0	2	0
			1576	1000	267	297	12			
1	B	187	Total	C	N	O	S	0	1	0
			1528	971	259	287	11			

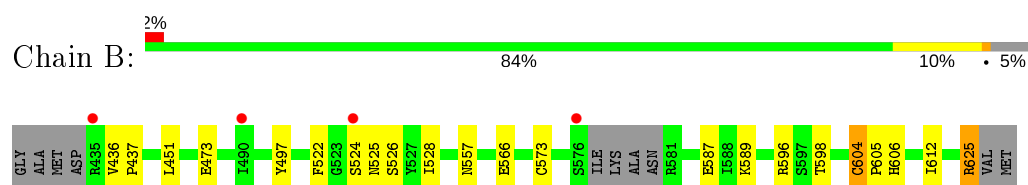
There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	431	GLY	-	EXPRESSION TAG	UNP P49850
A	432	ALA	-	EXPRESSION TAG	UNP P49850
A	433	MET	-	EXPRESSION TAG	UNP P49850
B	431	GLY	-	EXPRESSION TAG	UNP P49850
B	432	ALA	-	EXPRESSION TAG	UNP P49850
B	433	MET	-	EXPRESSION TAG	UNP P49850

- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	101	Total	O	0	0
			101	101		
2	B	89	Total	O	0	0
			89	89		

- Molecule 1: DNA mismatch repair protein mutL



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	33.42Å 74.33Å 182.47Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	38.88 – 2.00 38.88 – 1.99	Depositor EDS
% Data completeness (in resolution range)	99.2 (38.88-2.00) 99.2 (38.88-1.99)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	0.06	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.77 (at 1.98Å)	Xtriage
Refinement program	PHENIX (phenix.refine)	Depositor
R, R_{free}	0.191 , 0.228 0.191 , 0.226	Depositor DCC
R_{free} test set	1628 reflections (5.06%)	wwPDB-VP
Wilson B-factor (Å ²)	27.8	Xtriage
Anisotropy	0.132	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 45.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	3294	wwPDB-VP
Average B, all atoms (Å ²)	41.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 8.21% 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](#)

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	0.34	1/1615 (0.1%)	0.47	0/2175
1	B	0.37	1/1564 (0.1%)	0.47	0/2109
All	All	0.36	2/3179 (0.1%)	0.47	0/4284

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	587	GLU	CD-OE2	7.07	1.33	1.25
1	A	587	GLU	CD-OE2	6.20	1.32	1.25

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

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	1576	0	1551	33	0
1	B	1528	0	1501	18	0
2	A	101	0	0	3	0
2	B	89	0	0	1	0
All	All	3294	0	3052	46	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 7.

All (46) 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:608:ARG:HD3	1:B:625:ARG:NH1	1.99	0.77
1:A:604:CYS:SG	2:A:264:HOH:O	2.43	0.76
2:A:262:HOH:O	1:B:625:ARG:HG2	1.87	0.73
1:A:532:HIS:HD2	1:A:533:PRO:O	1.77	0.67
1:A:604:CYS:SG	1:A:608:ARG:HB2	2.34	0.67
1:A:622:MET:HE3	1:B:612:ILE:HG13	1.80	0.63
1:A:608:ARG:HD3	1:B:625:ARG:HH11	1.62	0.62
1:A:442:ILE:HG21	1:B:451:LEU:HD12	1.82	0.60
1:A:548:ILE:N	1:A:570:MET:HE1	2.17	0.59
1:A:512:LEU:HD21	1:A:552:VAL:HG21	1.84	0.58
1:B:497:TYR:O	1:B:525:ASN:HB3	2.02	0.58
1:B:573:CYS:SG	1:B:606:HIS:HE1	2.26	0.58
1:A:502:ALA:CB	1:A:525:ASN:HA	2.34	0.58
1:A:558:ILE:HD11	1:A:562:LYS:HG3	1.87	0.57
1:A:438:ILE:HA	1:A:596:ARG:HD2	1.86	0.57
1:A:548:ILE:HG13	1:A:570:MET:CE	2.35	0.56
1:B:473:GLU:HG3	1:B:598:THR:HG22	1.89	0.54
1:A:548:ILE:HG13	1:A:570:MET:HE2	1.90	0.52
1:B:522:PHE:CG	1:B:522:PHE:O	2.63	0.52
1:B:589:LYS:HE3	2:B:92:HOH:O	2.10	0.51
1:A:515:VAL:HG21	1:A:563:LEU:HG	1.93	0.51
1:A:489:MET:HE1	1:A:529:VAL:HG12	1.92	0.50
1:A:497:TYR:OH	1:A:549:ILE:HG21	2.12	0.50
1:A:544:LEU:HD21	1:A:571:MET:HG2	1.94	0.50
1:B:522:PHE:O	1:B:522:PHE:CD1	2.66	0.48
1:A:486:VAL:CG1	1:A:531:CYS:HB2	2.43	0.48
1:A:490:ILE:HG22	1:A:491:VAL:HG23	1.96	0.47
1:A:616:THR:HA	1:A:619:MET:HE2	1.95	0.47
1:A:515:VAL:O	1:A:515:VAL:CG1	2.62	0.47
1:B:436:VAL:HG22	1:B:437:PRO:O	2.15	0.47
1:B:522:PHE:CZ	1:B:526:SER:HB3	2.50	0.46
1:A:621:LYS:HA	1:A:627:MET:HE2	1.97	0.45
1:A:608:ARG:CD	1:B:625:ARG:NH1	2.76	0.45
1:B:604[A]:CYS:SG	1:B:605:PRO:HD2	2.57	0.45
1:B:473:GLU:HG3	1:B:598:THR:CG2	2.47	0.44
1:A:581:ARG:NH1	2:A:241:HOH:O	2.49	0.44
1:A:499:THR:HG22	1:A:500:ASN:OD1	2.19	0.43
1:B:436:VAL:HG13	1:B:596:ARG:HG3	2.01	0.42
1:A:619:MET:HA	1:A:622:MET:HE2	2.00	0.42
1:A:574:LYS:N	1:A:574:LYS:HD3	2.35	0.42
1:A:497:TYR:O	1:A:525:ASN:HB3	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:591:LEU:HD23	1:A:591:LEU:C	2.41	0.41
1:A:548:ILE:HG13	1:A:570:MET:HE1	2.03	0.41
1:A:515:VAL:O	1:A:515:VAL:HG12	2.20	0.41
1:A:508:HIS:HA	1:A:511:GLU:OE1	2.21	0.40
1:B:625:ARG:HD2	1:B:625:ARG:HA	1.81	0.40

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	189/197 (96%)	183 (97%)	6 (3%)	0	100	100
1	B	184/197 (93%)	181 (98%)	3 (2%)	0	100	100
All	All	373/394 (95%)	364 (98%)	9 (2%)	0	100	100

There are no Ramachandran outliers to report.

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	175/176 (99%)	167 (95%)	8 (5%)	27	23
1	B	169/176 (96%)	162 (96%)	7 (4%)	30	28
All	All	344/352 (98%)	329 (96%)	15 (4%)	30	25

All (15) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	488	GLU
1	A	499	THR
1	A	500	ASN
1	A	524	SER
1	A	553	LEU
1	A	562	LYS
1	A	608	ARG
1	A	612	ILE
1	B	524	SER
1	B	528	ILE
1	B	557	ASN
1	B	566	GLU
1	B	604[A]	CYS
1	B	604[B]	CYS
1	B	625	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	496	HIS
1	A	507	GLN
1	A	532	HIS
1	B	463	GLN
1	B	464	HIS
1	B	467	GLN
1	B	606	HIS

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

There are no ligands in this entry.

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(Å ²)	Q<0.9
1	A	191/197 (96%)	0.46	18 (9%) 8 8	14, 40, 96, 121	0
1	B	187/197 (94%)	-0.05	4 (2%) 63 62	13, 33, 64, 95	0
All	All	378/394 (95%)	0.21	22 (5%) 23 22	13, 36, 88, 121	0

All (22) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	522	PHE	5.6
1	A	553	LEU	5.5
1	B	576	SER	5.0
1	A	575	GLY	5.0
1	A	497	TYR	4.1
1	A	500	ASN	3.8
1	A	507	GLN	3.8
1	A	556	LYS	3.7
1	A	485	GLU	3.4
1	A	499	THR	3.3
1	B	435	ARG	2.9
1	A	490	ILE	2.7
1	A	558	ILE	2.7
1	B	524	SER	2.7
1	A	525	ASN	2.7
1	A	502	ALA	2.5
1	B	490	ILE	2.3
1	A	518	PHE	2.3
1	A	526	SER	2.2
1	A	557	ASN	2.1
1	A	486	VAL	2.0
1	A	549	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](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.