



wwPDB X-ray Structure Validation Summary Report ⓘ

Jan 31, 2016 – 09:05 PM GMT

PDB ID : 1NG9
Title : E.coli MutS R697A: an ATPase-asymmetry mutant
Authors : Lamers, M.H.; Winterwerp, H.H.K.; Sixma, T.K.
Deposited on : 2002-12-17
Resolution : 2.60 Å(reported)

This is a wwPDB X-ray Structure 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/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 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
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) : trunk26865

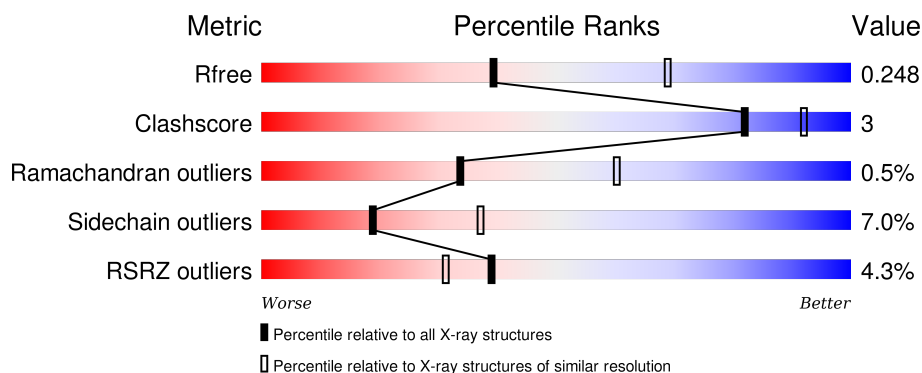
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.60 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}	91344	2328 (2.60-2.60)
Clashscore	102246	2679 (2.60-2.60)
Ramachandran outliers	100387	2635 (2.60-2.60)
Sidechain outliers	100360	2635 (2.60-2.60)
RSRZ outliers	91569	2334 (2.60-2.60)

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	E	30	<div> <div>13%</div> <div> <div></div> <div>43%</div> <div>13%</div> <div>•</div> <div>40%</div> </div> </div>
2	F	30	<div> <div>3%</div> <div> <div></div> <div>43%</div> <div>10%</div> <div>•</div> <div>43%</div> </div> </div>
3	A	800	<div> <div>4%</div> <div> <div></div> <div>85%</div> <div>13%</div> <div>••</div> </div> </div>
3	B	800	<div> <div>4%</div> <div> <div></div> <div>82%</div> <div>12%</div> <div>••</div> </div> </div>

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 13420 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 DNA chain called 5'-D(*AP*GP*CP*TP*GP*CP*CP*AP*GP*GP*CP*A
P*CP*CP*AP*GP*TP*GP*TP*CP*AP*GP*CP*GP*TP*CP*CP*TP*AP*T)-3'.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	E	18	Total	C	N	O	P	0	0	0
			367	174	72	104	17			

- Molecule 2 is a DNA chain called 5'-D(*AP*TP*AP*GP*GP*AP*CP*GP*CP*TP*GP*A
P*CP*AP*CP*TP*GP*GP*TP*GP*CP*TP*TP*GP*GP*CP*AP*GP*CP*T)-3'.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	F	17	Total	C	N	O	P	0	0	0
			347	166	62	103	16			

- Molecule 3 is a protein called DNA mismatch repair protein MutS.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	A	794	Total	C	N	O	S	0	0	0
			6241	3925	1109	1177	30			
3	B	766	Total	C	N	O	S	0	0	0
			6045	3808	1072	1136	29			

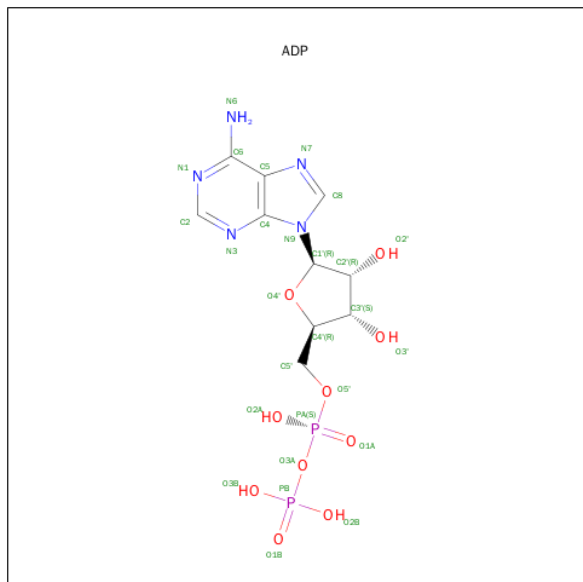
There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	697	ALA	ARG	ENGINEERED	UNP P23909
B	697	ALA	ARG	ENGINEERED	UNP P23909

- Molecule 4 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	B	1	Total	Mg	0	0
			1	1		
4	A	1	Total	Mg	0	0
			1	1		

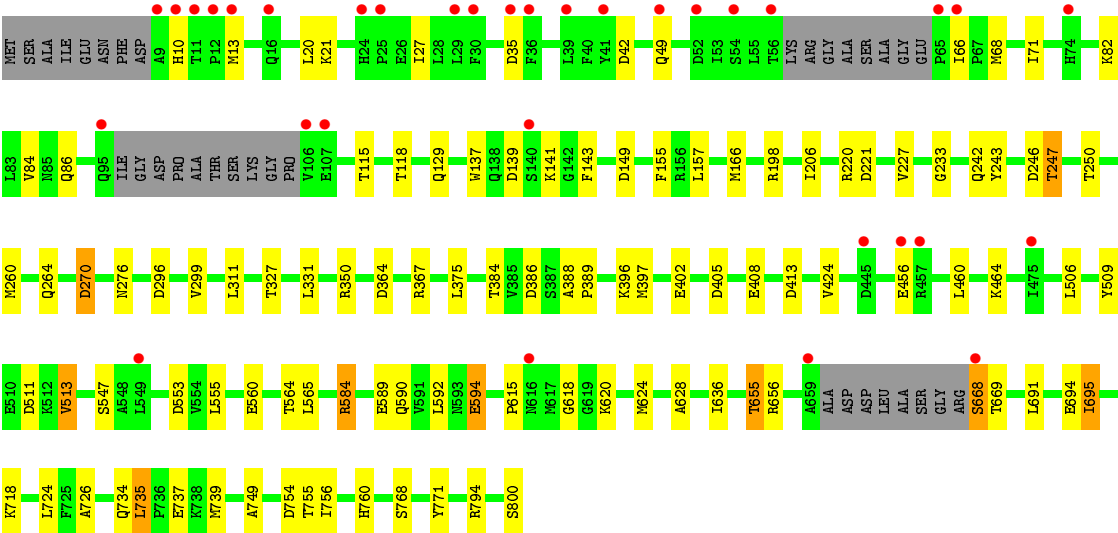
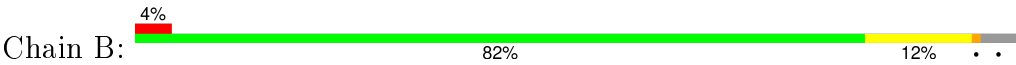
- Molecule 5 is ADENOSINE-5'-DIPHOSPHATE (three-letter code: ADP) (formula: $C_{10}H_{15}N_5O_{10}P_2$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
5	A	1	Total	C	N	O	P	0	0
			27	10	5	10	2		
5	B	1	Total	C	N	O	P	0	0
			27	10	5	10	2		

- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	165	Total	O	0	0
			165	165		
6	B	192	Total	O	0	0
			192	192		
6	E	3	Total	O	0	0
			3	3		
6	F	4	Total	O	0	0
			4	4		



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	89.90 Å 92.40 Å 261.59 Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.00 – 2.60 20.00 – 2.60	Depositor EDS
% Data completeness (in resolution range)	98.2 (20.00-2.60) 98.2 (20.00-2.60)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	0.12	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.09 (at 2.59 Å)	Xtriage
Refinement program	REFMAC 5.1.19	Depositor
R, R_{free}	0.215 , 0.249 0.212 , 0.248	Depositor DCC
R_{free} test set	1283 reflections (1.97%)	DCC
Wilson B-factor (Å ²)	39.1	Xtriage
Anisotropy	0.790	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 39.0	EDS
Estimated twinning fraction	0.024 for k,h,l	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Outliers	0 of 66556 reflections	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	13420	wwPDB-VP
Average B, all atoms (Å ²)	11.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.07% of the height of the origin peak. No significant pseudotranslation is detected.*

¹ Intensities estimated from amplitudes.

² Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.375 respectively for untwinned datasets, and 0.333, 0.2 for perfectly twinned datasets.

5 Model quality [i](#)

5.1 Standard geometry [i](#)

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

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	E	0.75	0/412	1.40	4/634 (0.6%)
2	F	0.71	0/388	1.50	4/598 (0.7%)
3	A	0.40	0/6347	0.73	12/8591 (0.1%)
3	B	0.42	0/6146	0.71	13/8318 (0.2%)
All	All	0.44	0/13293	0.79	33/18141 (0.2%)

There are no bond length outliers.

The worst 5 of 33 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	F	14	DA	O4'-C1'-N9	9.82	114.87	108.00
2	F	23	DT	O4'-C1'-N1	-7.59	102.69	108.00
1	E	1	DA	O4'-C1'-N9	7.51	113.26	108.00
2	F	18	DG	O4'-C4'-C3'	-7.22	101.61	104.50
3	B	270	ASP	CB-CG-OD2	6.55	124.20	118.30

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	E	367	0	202	1	0
2	F	347	0	194	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	A	6241	0	6288	43	0
3	B	6045	0	6095	38	0
4	A	1	0	0	0	0
4	B	1	0	0	0	0
5	A	27	0	12	0	0
5	B	27	0	12	1	0
6	A	165	0	0	9	0
6	B	192	0	0	5	0
6	E	3	0	0	1	0
6	F	4	0	0	0	0
All	All	13420	0	12803	82	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 3.

The worst 5 of 82 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)
6:E:358:HOH:O	3:A:58:ARG:HD3	1.85	0.76
3:B:327:THR:HG21	3:B:555:LEU:HD13	1.71	0.73
3:A:74:HIS:ND1	6:A:2120:HOH:O	2.20	0.72
3:A:74:HIS:CE1	6:A:2120:HOH:O	2.47	0.68
3:A:133:LEU:HG	3:A:240:LEU:HD11	1.78	0.64

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	
3	A	790/800 (99%)	765 (97%)	18 (2%)	7 (1%)	21	42
3	B	758/800 (95%)	733 (97%)	25 (3%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
All	All	1548/1600 (97%)	1498 (97%)	43 (3%)	7 (0%)	34 60

5 of 7 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
3	A	2	SER
3	A	594	GLU
3	A	439	LEU
3	A	3	ALA
3	A	140	SER

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
3	A	659/663 (99%)	615 (93%)	44 (7%)	20 40
3	B	640/663 (96%)	593 (93%)	47 (7%)	17 35
All	All	1299/1326 (98%)	1208 (93%)	91 (7%)	19 37

5 of 91 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
3	A	761	SER
3	B	71	ILE
3	B	718	LYS
3	A	768	SER
3	B	20	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 15 such sidechains are listed below:

Mol	Chain	Res	Type
3	B	49	GLN
3	B	214	ASN
3	B	370	HIS

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Mol	Chain	Res	Type
3	A	616	ASN
3	B	344	GLN

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 4 ligands modelled in this entry, 2 are monoatomic - leaving 2 for Mogul analysis.

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	$\# Z > 2$
5	ADP	A	2001	4	22,29,29	1.22	2 (9%)	27,45,45	2.72	4 (14%)
5	ADP	B	2002	4	22,29,29	1.14	2 (9%)	27,45,45	2.47	2 (7%)

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
5	ADP	A	2001	4	-	0/12/32/32	0/3/3/3
5	ADP	B	2002	4	-	0/12/32/32	0/3/3/3

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	B	2002	ADP	C2-N1	2.34	1.38	1.33
5	A	2001	ADP	C2-N1	2.87	1.39	1.33
5	B	2002	ADP	C2-N3	3.65	1.38	1.32
5	A	2001	ADP	C2-N3	3.76	1.38	1.32

The worst 5 of 6 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	A	2001	ADP	N3-C2-N1	-12.73	119.14	128.89
5	B	2002	ADP	N3-C2-N1	-11.36	120.19	128.89
5	B	2002	ADP	C2'-C1'-N9	-3.18	109.44	114.29
5	A	2001	ADP	PA-O3A-PB	-3.16	122.07	132.67
5	A	2001	ADP	C2'-C1'-N9	-2.01	111.23	114.29

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	B	2002	ADP	1	0

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 [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	E	18/30 (60%)	0.89	4 (22%) 1 0	4, 10, 16, 19	0
2	F	17/30 (56%)	0.37	1 (5%) 26 19	5, 10, 15, 18	0
3	A	794/800 (99%)	0.01	30 (3%) 44 36	4, 10, 14, 20	0
3	B	766/800 (95%)	0.07	33 (4%) 39 31	4, 10, 14, 23	0
All	All	1595/1660 (96%)	0.05	68 (4%) 39 31	4, 10, 14, 23	0

The worst 5 of 68 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
3	B	12	PRO	6.5
3	B	106	VAL	6.5
1	E	18	DG	6.4
3	B	659	ALA	6.4
1	E	1	DA	5.4

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
5	ADP	B	2002	27/27	0.90	0.20	0.20	64,71,73,76	0
5	ADP	A	2001	27/27	0.96	0.14	-0.95	50,56,58,58	0
4	MG	A	1001	1/1	0.89	0.05	-	41,41,41,41	0
4	MG	B	1002	1/1	0.90	0.45	-	54,54,54,54	0

6.5 Other polymers [i](#)

There are no such residues in this entry.