



## wwPDB EM Validation Summary Report ⓘ

May 12, 2024 – 12:07 AM JST

PDB ID : 8JQC  
EMDB ID : EMD-36569  
Title : Novel Anti-phage System  
Authors : Li, J.; Wang, Z.; Wang, L.  
Deposited on : 2023-06-13  
Resolution : 3.39 Å(reported)

This is a wwPDB EM Validation Summary 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/EMValidationReportHelp>

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>.

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

EMDB validation analysis : **FAILED**  
MolProbity : 4.02b-467  
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)  
MapQ : **FAILED**  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.36.2

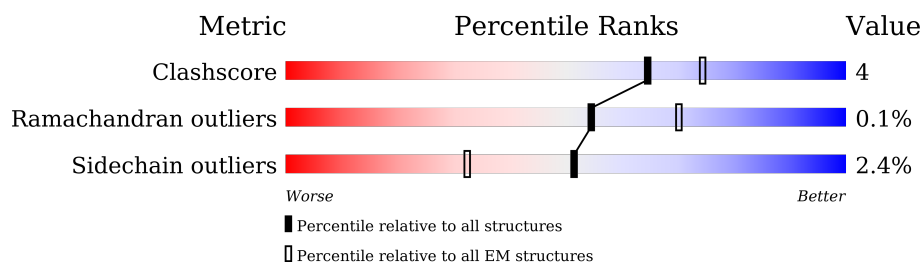
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*ELECTRON MICROSCOPY*

The reported resolution of this entry is 3.39 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)	EM structures (#Entries)
Clashscore	158937	4297
Ramachandran outliers	154571	4023
Sidechain outliers	154315	3826

The table below summarises the geometric issues observed across the polymeric chains and their fit to the map. The red, orange, yellow and green segments of the 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  $\leq 5\%$ .

Mol	Chain	Length	Quality of chain
1	A	578	70% 9% 21%
1	B	578	67% 12% 21%
1	C	578	68% 10% 21%
1	D	578	67% 11% 21%
2	E	499	81% 17% .

## 2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 18899 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, 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 Endonuclease GajA.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	454	Total	C	N	O	S	0	0
			3708	2378	613	705	12		
1	B	454	Total	C	N	O	S	0	0
			3708	2378	613	705	12		
1	C	454	Total	C	N	O	S	0	0
			3708	2378	613	705	12		
1	D	454	Total	C	N	O	S	0	0
			3708	2378	613	705	12		

- Molecule 2 is a protein called Gabija protein GajB.

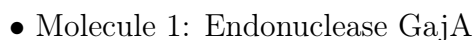
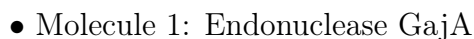
Mol	Chain	Residues	Atoms					AltConf	Trace
2	E	499	Total	C	N	O	S	0	0
			4067	2604	670	781	12		

There are 5 discrepancies between the modelled and reference sequences:

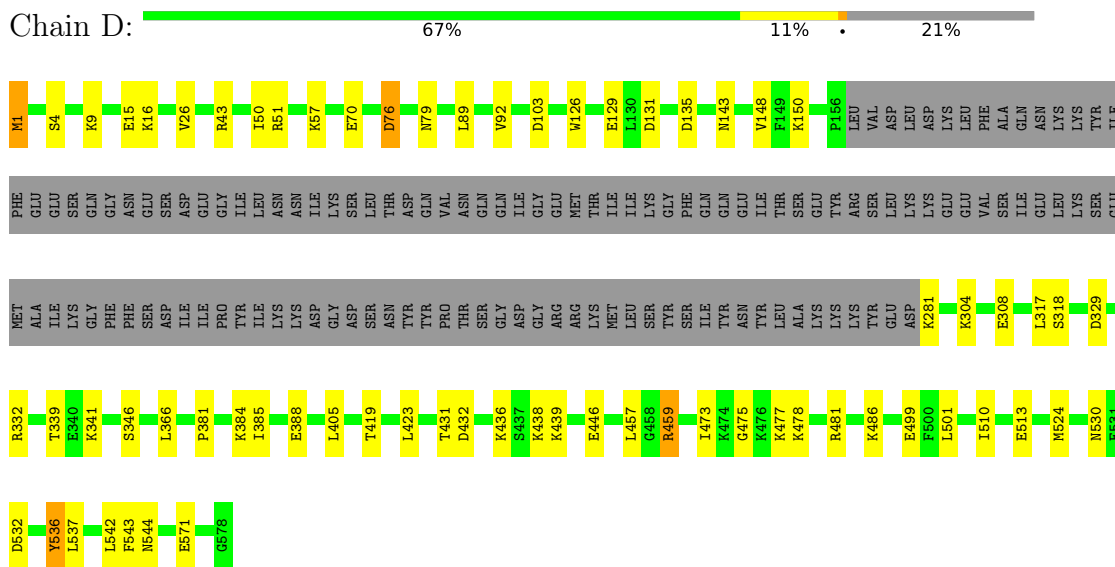
Chain	Residue	Modelled	Actual	Comment	Reference
E	1	MET	-	initiating methionine	UNP J8HQ06
E	2	ILE	-	expression tag	UNP J8HQ06
E	3	GLU	-	expression tag	UNP J8HQ06
E	4	ASP	-	expression tag	UNP J8HQ06
E	5	GLU	-	expression tag	UNP J8HQ06



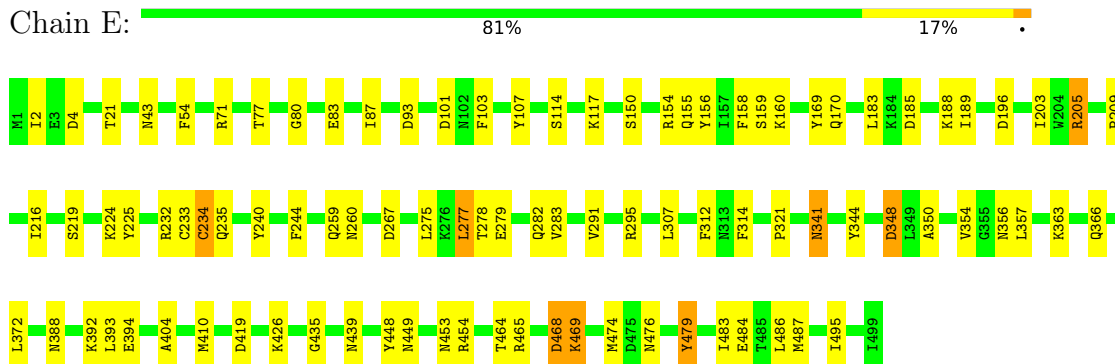
- Molecule 1: Endonuclease GajA



- Molecule 1: Endonuclease GajA



- Molecule 2: Gabija protein GajB



## 4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	4218276	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING AND AMPLITUDE CORRECTION	Depositor
Microscope	FEI TITAN KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ( $e^-/\text{\AA}^2$ )	50	Depositor
Minimum defocus (nm)	1600	Depositor
Maximum defocus (nm)	1800	Depositor
Magnification	Not provided	
Image detector	GATAN K3 (6k x 4k)	Depositor

## 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	0.27	0/3771	0.55	1/5064 (0.0%)
1	B	0.27	0/3771	0.57	3/5064 (0.1%)
1	C	0.28	0/3771	0.61	5/5064 (0.1%)
1	D	0.29	0/3771	0.60	2/5064 (0.0%)
2	E	0.31	0/4136	0.70	10/5579 (0.2%)
All	All	0.29	0/19220	0.61	21/25835 (0.1%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1
2	E	0	1
All	All	0	2

There are no bond length outliers.

The worst 5 of 21 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	E	209	PRO	CA-N-CD	-11.08	95.99	111.50
1	B	76	ASP	CB-CG-OD1	9.24	126.61	118.30
2	E	93	ASP	CB-CG-OD1	8.49	125.94	118.30
2	E	4	ASP	CB-CG-OD2	8.38	125.84	118.30
1	A	434	ASP	CB-CG-OD1	8.18	125.66	118.30

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	309	GLN	Peptide

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Mol	Chain	Res	Type	Group
2	E	267	ASP	Peptide

## 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	3708	0	3749	24	0
1	B	3708	0	3749	36	0
1	C	3708	0	3749	32	0
1	D	3708	0	3749	31	0
2	E	4067	0	4089	42	0
All	All	18899	0	19085	159	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.

The worst 5 of 159 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:304:LYS:O	1:A:308:GLU:HB3	1.96	0.66
1:D:332:ARG:HE	1:D:346:SER:HG	1.44	0.64
1:C:520:ILE:HD11	1:C:560:VAL:HB	1.82	0.62
2:E:307:LEU:HB3	2:E:312:PHE:HB2	1.83	0.61
1:B:96:ALA:HA	1:B:134:ILE:HD11	1.84	0.59

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 PDB entries followed by that with respect to all EM entries.

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	450/578 (78%)	434 (96%)	16 (4%)	0	100	100
1	B	450/578 (78%)	435 (97%)	15 (3%)	0	100	100
1	C	450/578 (78%)	428 (95%)	22 (5%)	0	100	100
1	D	450/578 (78%)	431 (96%)	19 (4%)	0	100	100
2	E	497/499 (100%)	437 (88%)	57 (12%)	3 (1%)	25	57
All	All	2297/2811 (82%)	2165 (94%)	129 (6%)	3 (0%)	54	82

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	E	354	VAL
2	E	2	ILE
2	E	203	ILE

### 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 PDB entries followed by that with respect to all EM entries.

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	416/529 (79%)	408 (98%)	8 (2%)	57	78
1	B	416/529 (79%)	410 (99%)	6 (1%)	67	83
1	C	416/529 (79%)	411 (99%)	5 (1%)	71	85
1	D	416/529 (79%)	401 (96%)	15 (4%)	35	63
2	E	453/453 (100%)	437 (96%)	16 (4%)	36	65
All	All	2117/2569 (82%)	2067 (98%)	50 (2%)	51	74

5 of 50 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	D	459	ARG
2	E	103	PHE
2	E	479	TYR

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Mol	Chain	Res	Type
1	D	499	GLU
1	D	543	PHE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

### 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 monosaccharides 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.