wwPDB X-ray Structure Validation Summary Report (i)

Dec 5, 2023 – 02:47 am GMT

PDB ID	:	1V14
Title	:	Crystal Structure of the Colicin E9, mutant His103Ala, in complex with Mg+2 and dsDNA (resolution 2.9A)
Authors	:	Mate, M.J.; Kleanthous, C.
Deposited on	:	2004-04-06
Resolution	:	2.90 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
https://www.wwpdb.org/validation/2017/XrayValidationReportHelp
with specific help available everywhere you see the (i) 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 (1)) were used in the production of this report:

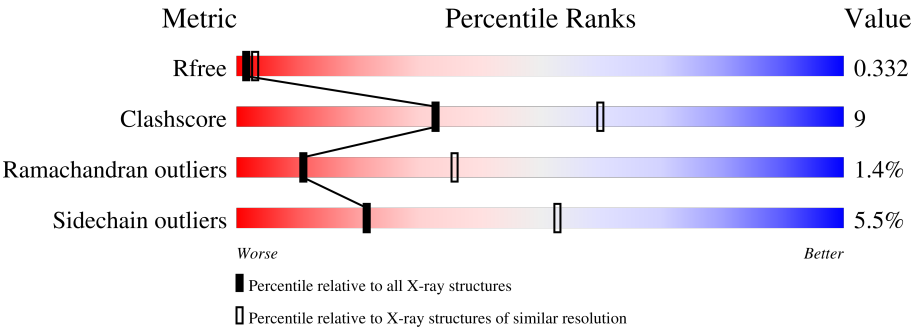
MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.36
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.90 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	1957 (2.90-2.90)
Clashscore	141614	2172 (2.90-2.90)
Ramachandran outliers	138981	2115 (2.90-2.90)
Sidechain outliers	138945	2117 (2.90-2.90)

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 >=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 <=5%

Mol	Chain	Length	Quality of chain
1	A	134	69% 27%
1	B	134	75% 20%
1	C	134	68% 27%
1	D	134	66% 30%
2	E	8	38% 38% 25%
2	F	8	38% 50% 12%
2	G	8	38% 38% 25%

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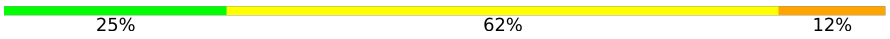
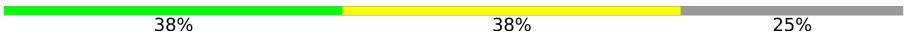
Mol	Chain	Length	Quality of chain
2	H	8	
2	I	8	
2	J	8	
2	K	8	
2	L	8	

2 Entry composition (i)

There are 4 unique types of molecules in this entry. The entry contains 5328 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 COLICIN E9.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	132	Total	C	N	O	S	0	0	0
			1041	647	192	201	1			
1	B	132	Total	C	N	O	S	0	0	0
			1045	650	192	201	2			
1	C	132	Total	C	N	O	S	0	0	0
			1041	647	192	201	1			
1	D	130	Total	C	N	O	S	0	0	0
			1026	639	190	196	1			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	103	ALA	HIS	engineered mutation	UNP P09883
B	103	ALA	HIS	engineered mutation	UNP P09883
C	103	ALA	HIS	engineered mutation	UNP P09883
D	103	ALA	HIS	engineered mutation	UNP P09883

- Molecule 2 is a DNA chain called 5'-D(*GP*CP*GP*AP*TP*CP*GP*CP)-3'.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	E	6	Total	C	N	O	P	0	0	0
			123	58	23	36	6			
2	F	8	Total	C	N	O	P	0	0	0
			161	77	31	46	7			
2	G	6	Total	C	N	O	P	0	0	0
			123	58	23	36	6			
2	H	8	Total	C	N	O	P	0	0	0
			161	77	31	46	7			
2	I	6	Total	C	N	O	P	0	0	0
			123	58	23	36	6			
2	J	8	Total	C	N	O	P	0	0	0
			161	77	31	46	7			

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	K	6	Total 123	C 58	N 23	O 36	P 6	0	0	0
2	L	8	Total 161	C 77	N 31	O 46	P 7	0	0	0

- Molecule 3 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total 1	Mg 1	0	0
3	C	1	Total 1	Mg 1	0	0
3	E	1	Total 1	Mg 1	0	0
3	G	1	Total 1	Mg 1	0	0
3	I	1	Total 1	Mg 1	0	0
3	K	1	Total 1	Mg 1	0	0

- Molecule 4 is water.

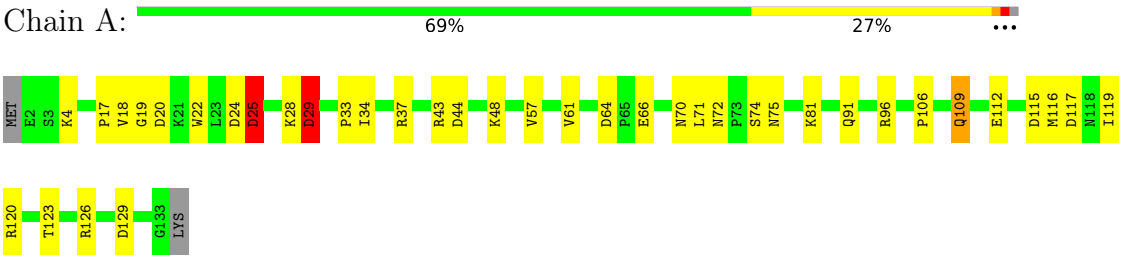
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	4	Total 4	O 4	0	0
4	B	5	Total 5	O 5	0	0
4	C	7	Total 7	O 7	0	0
4	D	6	Total 6	O 6	0	0
4	H	2	Total 2	O 2	0	0
4	I	1	Total 1	O 1	0	0
4	J	3	Total 3	O 3	0	0
4	K	3	Total 3	O 3	0	0
4	L	2	Total 2	O 2	0	0

3 Residue-property plots (i)

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. 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. 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.

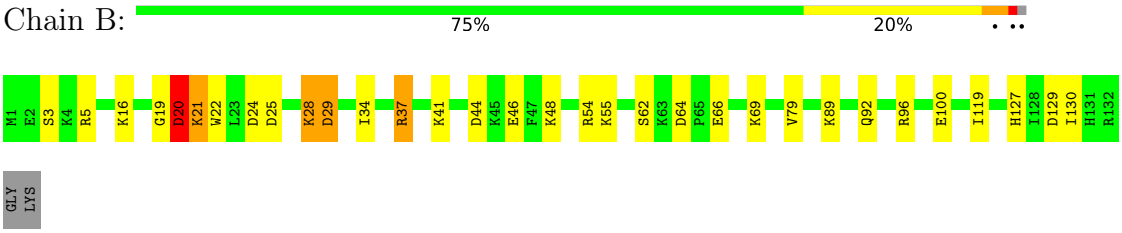
- Molecule 1: COLICIN E9

Chain A:



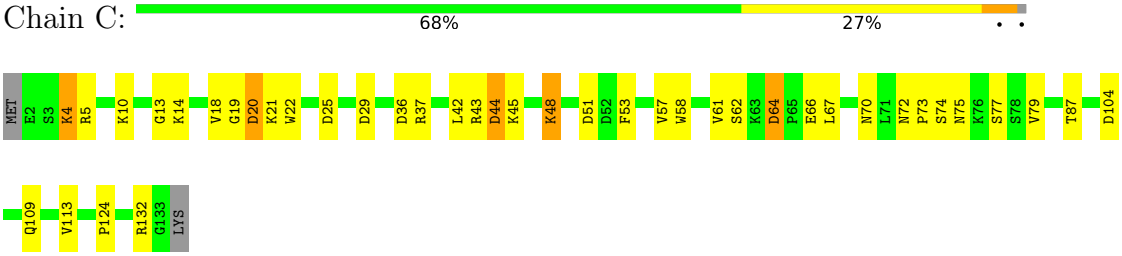
- Molecule 1: COLICIN E9

Chain B:



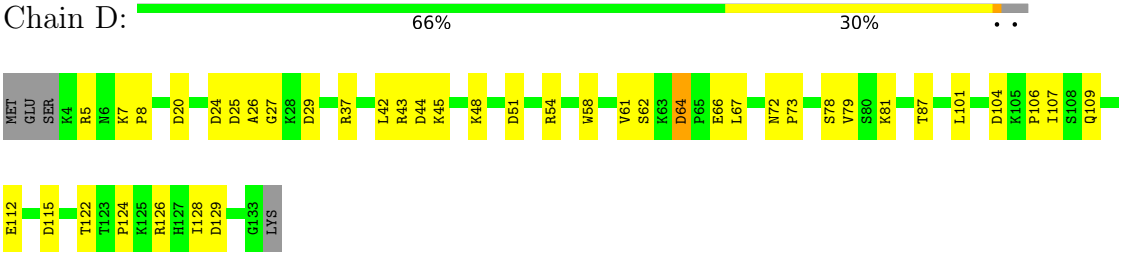
- Molecule 1: COLICIN E9

Chain C:



- Molecule 1: COLICIN E9

Chain D:



• Molecule 2: 5'-D(*GP*CP*GP*AP*TP*CP*GP*CP)-3'

Chain E:

• Molecule 2: 5'-D(*GP*CP*GP*AP*TP*CP*GP*CP)-3'

Chain F:

• Molecule 2: 5'-D(*GP*CP*GP*AP*TP*CP*GP*CP)-3'

Chain G:

• Molecule 2: 5'-D(*GP*CP*GP*AP*TP*CP*GP*CP)-3'

Chain H:

• Molecule 2: 5'-D(*GP*CP*GP*AP*TP*CP*GP*CP)-3'

Chain I:

• Molecule 2: 5'-D(*GP*CP*GP*AP*TP*CP*GP*CP)-3'

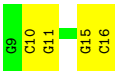
Chain J:

• Molecule 2: 5'-D(*GP*CP*GP*AP*TP*CP*GP*CP)-3'

Chain K:

• Molecule 2: 5'-D(*GP*CP*GP*AP*TP*CP*GP*CP)-3'

Chain L:


G9
C10
G11
G15
C16

4 Data and refinement statistics (i)

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α, β, γ	92.95Å 124.44Å 111.23Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	74.54 – 2.90 55.61 – 2.90	Depositor EDS
% Data completeness (in resolution range)	99.7 (74.54-2.90) 99.7 (55.61-2.90)	Depositor EDS
R_{merge}	0.04	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	3.26 (at 2.91Å)	Xtriage
Refinement program	REFMAC 5.2.0001	Depositor
R, R_{free}	0.215 , 0.306 0.251 , 0.332	Depositor DCC
R_{free} test set	734 reflections (5.03%)	wwPDB-VP
Wilson B-factor (Å^2)	52.3	Xtriage
Anisotropy	0.502	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.30 , 72.0	EDS
L-test for twinning[2]	$< \|L\| > = 0.46$, $< L^2 > = 0.29$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.91	EDS
Total number of atoms	5328	wwPDB-VP
Average B, all atoms (Å^2)	56.0	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 52.05 % 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 5.1415e-05. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ 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: MG

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	0.40	0/1064	0.81	7/1428 (0.5%)
1	B	0.51	1/1068 (0.1%)	0.82	7/1433 (0.5%)
1	C	0.43	0/1064	0.84	7/1428 (0.5%)
1	D	0.43	0/1049	0.83	9/1408 (0.6%)
2	E	0.76	0/137	1.56	4/209 (1.9%)
2	F	0.84	0/180	1.83	7/276 (2.5%)
2	G	0.86	0/137	1.46	3/209 (1.4%)
2	H	0.96	0/180	2.16	12/276 (4.3%)
2	I	0.78	0/137	1.53	1/209 (0.5%)
2	J	0.96	0/180	1.97	7/276 (2.5%)
2	K	0.76	0/137	1.43	1/209 (0.5%)
2	L	0.86	0/180	1.59	2/276 (0.7%)
All	All	0.57	1/5513 (0.0%)	1.13	67/7637 (0.9%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	16	LYS	CE-NZ	5.94	1.64	1.49

The worst 5 of 67 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	H	10	DC	O4'-C1'-N1	9.90	114.93	108.00
2	J	15	DG	OP2-P-O3'	-9.76	83.72	105.20
2	J	10	DC	O4'-C1'-N1	9.75	114.82	108.00
2	E	5	DT	C6-C5-C7	-9.19	117.39	122.90
2	F	15	DG	OP2-P-O3'	-9.17	85.02	105.20

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	1041	0	1033	17	0
1	B	1045	0	1042	13	0
1	C	1041	0	1033	27	0
1	D	1026	0	1022	21	0
2	E	123	0	68	0	0
2	F	161	0	91	3	0
2	G	123	0	68	2	0
2	H	161	0	91	2	0
2	I	123	0	68	1	0
2	J	161	0	91	3	0
2	K	123	0	68	0	0
2	L	161	0	91	1	0
3	A	1	0	0	0	0
3	C	1	0	0	0	0
3	E	1	0	0	0	0
3	G	1	0	0	0	0
3	I	1	0	0	0	0
3	K	1	0	0	0	0
4	A	4	0	0	0	0
4	B	5	0	0	0	0
4	C	7	0	0	0	0
4	D	6	0	0	0	0
4	H	2	0	0	0	0
4	I	1	0	0	0	0
4	J	3	0	0	0	0
4	K	3	0	0	0	0
4	L	2	0	0	0	0
All	All	5328	0	4766	87	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 9.

The worst 5 of 87 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:29:ASP:HA	1:A:126:ARG:HH21	1.40	0.87
2:F:10:DC:H2"	2:F:11:DG:O5'	1.85	0.76
1:A:34:ILE:HG21	1:A:116:MET:O	1.92	0.70
1:D:54:ARG:HG3	1:D:101:LEU:HD12	1.72	0.69
1:C:14:LYS:HA	1:C:43:ARG:NH1	2.08	0.68

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	130/134 (97%)	116 (89%)	13 (10%)	1 (1%)	19 51
1	B	130/134 (97%)	115 (88%)	11 (8%)	4 (3%)	4 16
1	C	130/134 (97%)	117 (90%)	11 (8%)	2 (2%)	10 34
1	D	128/134 (96%)	120 (94%)	8 (6%)	0	100 100
All	All	518/536 (97%)	468 (90%)	43 (8%)	7 (1%)	11 36

5 of 7 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	20	ASP
1	A	29	ASP
1	B	21	LYS
1	C	132	ARG
1	B	3	SER

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	114/116 (98%)	104 (91%)	10 (9%)	10 30
1	B	115/116 (99%)	108 (94%)	7 (6%)	18 48
1	C	114/116 (98%)	109 (96%)	5 (4%)	28 61
1	D	112/116 (97%)	109 (97%)	3 (3%)	44 77
All	All	455/464 (98%)	430 (94%)	25 (6%)	21 53

5 of 25 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	48	LYS
1	C	5	ARG
1	D	81	LYS
1	B	69	LYS
1	C	20	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	75	ASN
1	B	102	HIS
1	C	6	ASN
1	D	102	HIS
1	D	127	HIS

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)

Of 6 ligands modelled in this entry, 6 are monoatomic - leaving 0 for Mogul analysis.

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 (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

Unable to reproduce the depositors R factor - this section is therefore empty.

6.2 Non-standard residues in protein, DNA, RNA chains

Unable to reproduce the depositors R factor - this section is therefore empty.

6.3 Carbohydrates

Unable to reproduce the depositors R factor - this section is therefore empty.

6.4 Ligands

Unable to reproduce the depositors R factor - this section is therefore empty.

6.5 Other polymers

Unable to reproduce the depositors R factor - this section is therefore empty.