



Full wwPDB X-ray Structure Validation Report ⓘ

Jan 30, 2021 – 07:25 PM EST

PDB ID : 3LYB
Title : Structure of putative endoribonuclease(KP1_3112) from *Klebsiella pneumoniae*
Authors : Ramagopal, U.A.; Toro, R.; Burley, S.K.; Almo, S.C.; New York SGX Research Center for Structural Genomics (NYSGXRC)
Deposited on : 2010-02-26
Resolution : 2.66 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
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.16
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) : 2.16

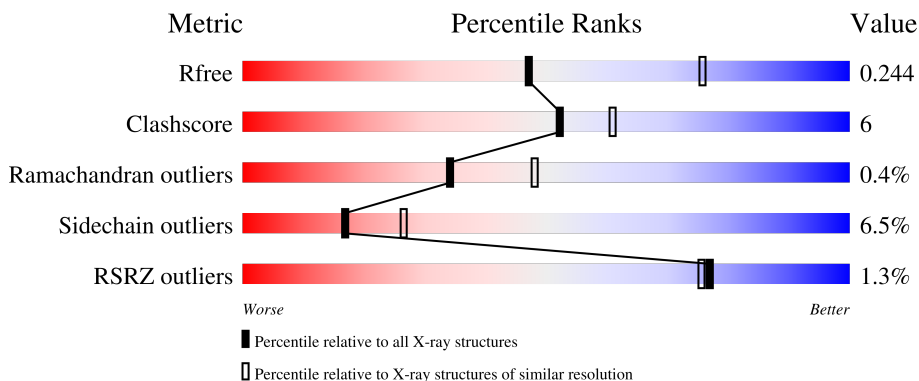
1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.66 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	1332 (2.68-2.64)
Clashscore	141614	1374 (2.68-2.64)
Ramachandran outliers	138981	1349 (2.68-2.64)
Sidechain outliers	138945	1349 (2.68-2.64)
RSRZ outliers	127900	1318 (2.68-2.64)

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 $\leq 5\%$. 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	165	<div style="display: flex; align-items: center;"> <div style="width: 2%; height: 10px; background-color: red; margin-right: 2px;"></div> <div style="width: 72%; height: 10px; background-color: green; margin-right: 2px;"></div> <div style="width: 9%; height: 10px; background-color: yellow; margin-right: 2px;"></div> <div style="width: 1%; height: 10px; background-color: orange; margin-right: 2px;"></div> <div style="width: 18%; height: 10px; background-color: grey; margin-right: 2px;"></div> </div> <p style="margin-left: 20px;">72% 9% • 18%</p>
1	B	165	<div style="display: flex; align-items: center;"> <div style="width: 2%; height: 10px; background-color: red; margin-right: 2px;"></div> <div style="width: 67%; height: 10px; background-color: green; margin-right: 2px;"></div> <div style="width: 15%; height: 10px; background-color: yellow; margin-right: 2px;"></div> <div style="width: 1%; height: 10px; background-color: orange; margin-right: 2px;"></div> <div style="width: 17%; height: 10px; background-color: grey; margin-right: 2px;"></div> </div> <p style="margin-left: 20px;">67% 15% • 17%</p>
1	C	165	<div style="display: flex; align-items: center;"> <div style="width: 2%; height: 10px; background-color: red; margin-right: 2px;"></div> <div style="width: 65%; height: 10px; background-color: green; margin-right: 2px;"></div> <div style="width: 14%; height: 10px; background-color: yellow; margin-right: 2px;"></div> <div style="width: 1%; height: 10px; background-color: orange; margin-right: 2px;"></div> <div style="width: 19%; height: 10px; background-color: grey; margin-right: 2px;"></div> </div> <p style="margin-left: 20px;">65% 14% • 19%</p>
1	D	165	<div style="display: flex; align-items: center;"> <div style="width: 2%; height: 10px; background-color: red; margin-right: 2px;"></div> <div style="width: 64%; height: 10px; background-color: green; margin-right: 2px;"></div> <div style="width: 16%; height: 10px; background-color: yellow; margin-right: 2px;"></div> <div style="width: 1%; height: 10px; background-color: orange; margin-right: 2px;"></div> <div style="width: 18%; height: 10px; background-color: grey; margin-right: 2px;"></div> </div> <p style="margin-left: 20px;">64% 16% • 18%</p>

2 Entry composition i

There are 3 unique types of molecules in this entry. The entry contains 4290 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 Putative endoribonuclease.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	Se			
1	A	136	1071	695	178	196	2	0	0	0
1	B	137	1071	694	179	196	2	0	0	0
1	C	134	1058	686	176	194	2	0	0	0
1	D	135	1058	688	175	193	2	0	0	0

There are 32 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	158	GLU	-	expression tag	UNP C4X9E0
A	159	GLY	-	expression tag	UNP C4X9E0
A	160	HIS	-	expression tag	UNP C4X9E0
A	161	HIS	-	expression tag	UNP C4X9E0
A	162	HIS	-	expression tag	UNP C4X9E0
A	163	HIS	-	expression tag	UNP C4X9E0
A	164	HIS	-	expression tag	UNP C4X9E0
A	165	HIS	-	expression tag	UNP C4X9E0
B	158	GLU	-	expression tag	UNP C4X9E0
B	159	GLY	-	expression tag	UNP C4X9E0
B	160	HIS	-	expression tag	UNP C4X9E0
B	161	HIS	-	expression tag	UNP C4X9E0
B	162	HIS	-	expression tag	UNP C4X9E0
B	163	HIS	-	expression tag	UNP C4X9E0
B	164	HIS	-	expression tag	UNP C4X9E0
B	165	HIS	-	expression tag	UNP C4X9E0
C	158	GLU	-	expression tag	UNP C4X9E0
C	159	GLY	-	expression tag	UNP C4X9E0
C	160	HIS	-	expression tag	UNP C4X9E0
C	161	HIS	-	expression tag	UNP C4X9E0
C	162	HIS	-	expression tag	UNP C4X9E0

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Chain	Residue	Modelled	Actual	Comment	Reference
C	163	HIS	-	expression tag	UNP C4X9E0
C	164	HIS	-	expression tag	UNP C4X9E0
C	165	HIS	-	expression tag	UNP C4X9E0
D	158	GLU	-	expression tag	UNP C4X9E0
D	159	GLY	-	expression tag	UNP C4X9E0
D	160	HIS	-	expression tag	UNP C4X9E0
D	161	HIS	-	expression tag	UNP C4X9E0
D	162	HIS	-	expression tag	UNP C4X9E0
D	163	HIS	-	expression tag	UNP C4X9E0
D	164	HIS	-	expression tag	UNP C4X9E0
D	165	HIS	-	expression tag	UNP C4X9E0

- Molecule 2 is CALCIUM ION (three-letter code: CA) (formula: Ca).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	B	2	Total Ca 2 2	0	0
2	A	1	Total Ca 1 1	0	0
2	D	1	Total Ca 1 1	0	0

- Molecule 3 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	10	Total O 10 10	0	0
3	B	10	Total O 10 10	0	0
3	C	1	Total O 1 1	0	0
3	D	7	Total O 7 7	0	0

V124	S125	Q129	M133	L134	P135	P150	PRO	SER	PHE	ASN	SER	GLU	ALA	GLU	GLY	HIS	HIS	HIS	HIS	HIS	HIS	HIS
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4 Data and refinement statistics

Property	Value	Source
Space group	I 2 3	Depositor
Cell constants a, b, c, α , β , γ	166.56Å 166.56Å 166.56Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	37.24 – 2.66 35.51 – 2.66	Depositor EDS
% Data completeness (in resolution range)	99.8 (37.24-2.66) 99.8 (35.51-2.66)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	5.62 (at 2.65Å)	Xtrriage
Refinement program	REFMAC	Depositor
R, R_{free}	0.193 , 0.248 0.193 , 0.244	Depositor DCC
R_{free} test set	1122 reflections (5.05%)	wwPDB-VP
Wilson B-factor (Å ²)	47.4	Xtrriage
Anisotropy	0.000	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 60.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.44$, $\langle L^2 \rangle = 0.26$	Xtrriage
Estimated twinning fraction	0.055 for -l,-k,-h	Xtrriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	4290	wwPDB-VP
Average B, all atoms (Å ²)	40.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.44% 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](#)

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

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	0.70	0/1098	0.73	0/1499
1	B	0.65	1/1098 (0.1%)	0.70	0/1499
1	C	0.64	0/1084	0.69	0/1478
1	D	0.65	0/1084	0.70	0/1478
All	All	0.66	1/4364 (0.0%)	0.71	0/5954

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	52	GLU	CB-CG	5.12	1.61	1.52

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	1071	0	1068	7	0
1	B	1071	0	1056	18	0
1	C	1058	0	1051	12	0
1	D	1058	0	1041	12	0
2	A	1	0	0	0	0
2	B	2	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	D	1	0	0	0	0
3	A	10	0	0	0	0
3	B	10	0	0	1	0
3	C	1	0	0	0	0
3	D	7	0	0	0	0
All	All	4290	0	4216	47	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 6.

All (47) 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:B:121:GLN:HE21	1:B:121:GLN:H	1.34	0.74
1:C:104:LEU:HG	1:C:129:GLN:HB2	1.81	0.62
1:A:91:MSE:HE3	1:A:118:TYR:CZ	2.35	0.61
1:A:133:MSE:HE2	1:A:139:VAL:HG12	1.84	0.60
1:A:62:THR:O	1:A:66:GLN:HG2	2.02	0.59
1:B:30:ILE:HD12	1:B:32:VAL:HG13	1.85	0.59
1:B:81:ARG:HH11	1:B:81:ARG:HG2	1.68	0.58
1:C:32:VAL:HG12	1:C:39:ILE:HG13	1.86	0.57
1:B:104:LEU:HG	1:B:129:GLN:HB2	1.88	0.54
1:D:104:LEU:HG	1:D:129:GLN:HB2	1.89	0.54
1:D:81:ARG:HG3	1:D:91:MSE:HG2	1.89	0.54
1:B:36:THR:CB	1:B:38:THR:OG1	2.56	0.54
1:B:24:ILE:HB	1:B:147:VAL:HB	1.90	0.53
1:D:33:ASN:HB2	1:D:36:THR:O	2.09	0.52
1:B:30:ILE:HD12	1:B:32:VAL:CG1	2.41	0.50
1:D:62:THR:O	1:D:66:GLN:HG2	2.12	0.50
1:B:118:TYR:CZ	1:B:123:PRO:HG3	2.48	0.49
1:D:48:GLU:HG3	1:D:49:PRO:HD3	1.95	0.49
1:B:121:GLN:O	1:B:121:GLN:HG2	2.14	0.48
1:C:68:PRO:O	1:C:72:GLN:HG3	2.13	0.48
1:A:113:VAL:O	1:A:116:LEU:HB2	2.14	0.47
1:A:46:VAL:HG23	1:A:51:ARG:HG3	1.96	0.47
1:B:62:THR:O	1:B:66:GLN:HG2	2.14	0.47
1:C:20:ALA:HB3	1:C:25:PHE:HE1	1.80	0.46
1:D:33:ASN:N	1:D:33:ASN:OD1	2.49	0.46
1:B:121:GLN:N	1:B:121:GLN:HE21	2.08	0.46
1:B:82:ARG:NH1	3:B:171:HOH:O	2.49	0.46
1:C:24:ILE:HB	1:C:147:VAL:HB	1.98	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:107:PHE:HB3	1:B:108:PRO:HD3	1.98	0.45
1:A:59:GLU:OE2	1:C:115:LYS:HE2	2.17	0.44
1:D:31:PRO:HB2	1:D:40:VAL:HB	1.99	0.44
1:C:31:PRO:HD2	1:C:139:VAL:HG13	2.00	0.44
1:B:121:GLN:HA	1:B:122:PRO:HD2	1.87	0.43
1:B:69:ILE:HG12	1:B:106:HIS:HB3	2.01	0.43
1:B:60:PHE:CE2	1:C:135:PRO:HG3	2.54	0.43
1:C:29:ILE:HD12	1:C:80:ILE:HG12	2.01	0.42
1:D:69:ILE:HG12	1:D:106:HIS:HB3	2.02	0.42
1:B:78:GLU:OE2	1:B:81:ARG:NH2	2.53	0.42
1:D:43:PHE:O	1:D:46:VAL:HG22	2.19	0.42
1:D:104:LEU:HD23	1:D:104:LEU:HA	1.89	0.41
1:C:40:VAL:HG21	1:C:75:TYR:CD1	2.55	0.41
1:A:95:ILE:HG13	1:A:95:ILE:H	1.46	0.41
1:B:81:ARG:NH1	1:B:81:ARG:HG2	2.34	0.41
1:D:39:ILE:HD11	1:D:135:PRO:HG2	2.03	0.41
1:C:57:THR:HG21	1:C:62:THR:HB	2.02	0.40
1:C:36:THR:HB	1:C:38:THR:OG1	2.21	0.40
1:D:22:ASP:HB3	1:D:150:PRO:HG3	2.03	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	134/165 (81%)	129 (96%)	5 (4%)	0	100	100
1	B	135/165 (82%)	128 (95%)	6 (4%)	1 (1%)	22	33
1	C	130/165 (79%)	125 (96%)	5 (4%)	0	100	100
1	D	131/165 (79%)	123 (94%)	7 (5%)	1 (1%)	19	29
All	All	530/660 (80%)	505 (95%)	23 (4%)	2 (0%)	34	48

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	120	ASP
1	D	120	ASP

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	117/135 (87%)	110 (94%)	7 (6%)	19	30
1	B	115/135 (85%)	111 (96%)	4 (4%)	36	52
1	C	116/135 (86%)	109 (94%)	7 (6%)	19	30
1	D	113/135 (84%)	101 (89%)	12 (11%)	6	10
All	All	461/540 (85%)	431 (94%)	30 (6%)	17	26

All (30) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	32	VAL
1	A	41	ASN
1	A	95	ILE
1	A	99	GLN
1	A	111	SER
1	A	116	LEU
1	A	120	ASP
1	B	38	THR
1	B	92	SER
1	B	99	GLN
1	B	121	GLN
1	C	18	ARG
1	C	33	ASN
1	C	36	THR
1	C	38	THR
1	C	52	GLU
1	C	57	THR
1	C	96	LYS

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Mol	Chain	Res	Type
1	D	19	ARG
1	D	33	ASN
1	D	36	THR
1	D	48	GLU
1	D	52	GLU
1	D	86	SER
1	D	95	ILE
1	D	96	LYS
1	D	99	GLN
1	D	124	VAL
1	D	125	SER
1	D	133	MSE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	90	GLN
1	B	121	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 monosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 4 ligands modelled in this entry, 4 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 [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	134/165 (81%)	-0.26	1 (0%) 87 87	23, 36, 56, 63	0
1	B	135/165 (81%)	-0.23	1 (0%) 87 87	24, 39, 62, 69	0
1	C	132/165 (80%)	-0.03	3 (2%) 60 56	23, 42, 59, 62	0
1	D	133/165 (80%)	-0.14	2 (1%) 73 71	25, 43, 58, 63	0
All	All	534/660 (80%)	-0.17	7 (1%) 77 75	23, 40, 60, 69	0

All (7) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	35	LEU	3.4
1	D	36	THR	3.4
1	C	36	THR	3.0
1	C	34	PRO	2.4
1	A	35	LEU	2.4
1	C	86	SER	2.3
1	D	48	GLU	2.1

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 monosaccharides 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. 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	B-factors(Å ²)	Q<0.9
2	CA	D	166	1/1	0.89	0.04	72,72,72,72	0
2	CA	B	166	1/1	0.90	0.13	52,52,52,52	1
2	CA	A	166	1/1	0.93	0.06	79,79,79,79	0
2	CA	B	167	1/1	0.98	0.09	72,72,72,72	0

6.5 Other polymers [i](#)

There are no such residues in this entry.