

wwPDB X-ray Structure Validation Summary Report (i)

Feb 11, 2024 – 04:45 PM EST

PDB ID : 3DIL

Title : Crystal structure of the Thermotoga maritima lysine riboswitch bound to ly-

sine

Authors : Serganov, A.A. Deposited on : 2008-06-20

Resolution : 1.90 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:

Mol Probity : 4.02b-467

Mogul: 1.8.5 (274361), CSD as541be (2020)

Xtriage (Phenix) : 1.13 EDS : 2.36

Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)

 $Refmac \quad : \quad 5.8.0158$

CCP4 : 7.0.044 (Gargrove)

Ideal geometry (proteins) : Engh & Huber (2001) Ideal geometry (DNA, RNA) : Parkinson et al. (1996)

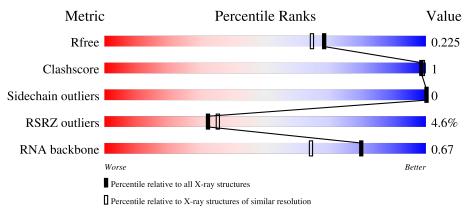
Validation Pipeline (wwPDB-VP) : 2.36

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 1.90 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 $(\# \mathrm{Entries})$	$\begin{array}{c} {\rm Similar\ resolution} \\ (\#{\rm Entries},\ {\rm resolution\ range}({\rm \AA})) \end{array}$
R_{free}	130704	6207 (1.90-1.90)
Clashscore	141614	6847 (1.90-1.90)
Sidechain outliers	138945	6760 (1.90-1.90)
RSRZ outliers	127900	6082 (1.90-1.90)
RNA backbone	3102	1013 (2.42-1.38)

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% 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			
ſ				5%			
	1	A	174	83%	16%	•	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
6	1PE	A	250	-	-	-	X



2 Entry composition (i)

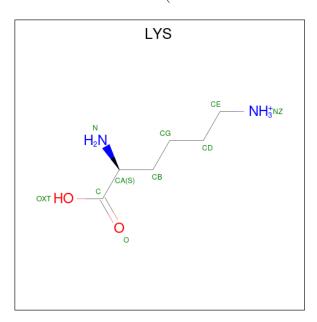
There are 8 unique types of molecules in this entry. The entry contains 4603 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 RNA chain called RNA (174-MER).

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace		
1	A	174	Total 3752	C 1666	N 695	O 1216	P 175	0	0	0

• Molecule 2 is LYSINE (three-letter code: LYS) (formula: $C_6H_{15}N_2O_2$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total 10	C 6	N 2	O 2	0	0

• Molecule 3 is POTASSIUM ION (three-letter code: K) (formula: K).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	3	Total K 3 3	0	0

• Molecule 4 is SODIUM ION (three-letter code: NA) (formula: Na).

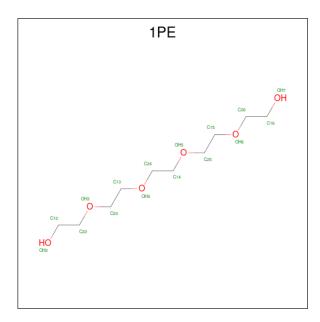


Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	29	Total Na 29 29	0	0

• Molecule 5 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	1	Total Mg 1 1	0	0

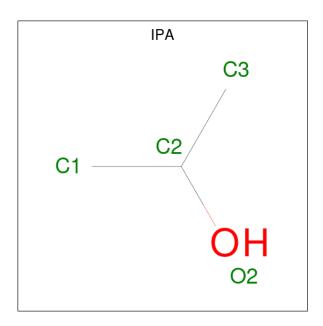
• Molecule 6 is PENTAETHYLENE GLYCOL (three-letter code: 1PE) (formula: $C_{10}H_{22}O_6$).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	A	1	Total C O 3 2 1	0	0
6	A	1	Total C O 3 2 1	0	0
6	A	1	Total C O 5 4 1	0	0

• Molecule 7 is ISOPROPYL ALCOHOL (three-letter code: IPA) (formula: C₃H₈O).





Mol	Chain	Residues	Atoms		ZeroOcc	AltConf	
7	A	1	Total 4	C 3	O 1	0	0

• Molecule 8 is water.

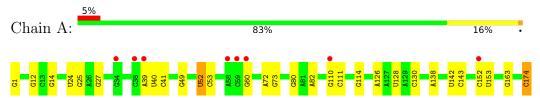
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
8	A	793	Total O 793 793	0	0



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 and electron density. 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. A red dot above a residue indicates a poor fit to the electron density (RSRZ > 2). 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: RNA (174-MER)





4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants	54.22Å 78.95Å 139.99Å	Donositor
a, b, c, α , β , γ	90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.00 - 1.90	Depositor
Resolution (A)	19.74 - 1.90	EDS
% Data completeness	95.8 (20.00-1.90)	Depositor
(in resolution range)	95.8 (19.74-1.90)	EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) > 1$	3.14 (at 1.90Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
D D.	0.192 , 0.229	Depositor
R, R_{free}	0.189 , 0.225	DCC
R_{free} test set	2334 reflections (5.08%)	wwPDB-VP
Wilson B-factor (Å ²)	20.6	Xtriage
Anisotropy	0.162	Xtriage
Bulk solvent $k_{sol}(e/Å^3)$, $B_{sol}(Å^2)$	0.35, 41.2	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	0.94	EDS
Total number of atoms	4603	wwPDB-VP
Average B, all atoms (Å ²)	23.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: The largest off-origin peak in the Patterson function is 5.89% of the height of the origin peak. No significant pseudotranslation is detected.

²Theoretical 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.



¹Intensities estimated from amplitudes.

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, CCC, K, 1PE, IPA, NA

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	Boı	nd lengths	Bo	nd angles
MOI	Chain	RMSZ	# Z > 5	RMSZ	# Z > 5
1	A	0.62	1/4174 (0.0%)	1.07	6/6514 (0.1%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	\mathbf{Z}	$\operatorname{Observed}(\text{\AA})$	$Ideal(\AA)$
1	A	1	G	OP3-P	-10.35	1.48	1.61

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

Mol	Chain	Res	Type	Atoms	Z	$\mathbf{Observed}(^{o})$	$\operatorname{Ideal}({}^{o})$
1	A	12	G	O4'-C1'-N9	8.91	115.33	108.20
1	A	52	U	P-O3'-C3'	6.07	126.99	119.70
1	A	27	G	O4'-C1'-N9	5.94	112.95	108.20
1	A	24	U	O4'-C1'-N1	5.80	112.84	108.20
1	A	130	С	O4'-C1'-N1	5.44	112.55	108.20

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	3752	0	1884	3	0	
2	A	10	0	12	1	0	

Continued on next page...



Continued	trom	mmoninonic	maaa
COHABABACA		DIEUIUU	DUIUE
0 0 1000100000			

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	A	3	0	0	0	0
4	A	29	0	0	0	0
5	A	1	0	0	0	0
6	A	11	0	10	0	0
7	A	4	0	8	0	0
8	A	793	0	0	0	0
All	All	4603	0	1914	3	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 1.

All (3) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	$\begin{array}{c} \text{Interatomic} \\ \text{distance (Å)} \end{array}$	Clash overlap (Å)
1:A:114:G:HO2'	2:A:175:LYS:N	1.97	0.61
1:A:25:G:H4'	1:A:126:A:C4	2.50	0.47
1:A:142:U:H2'	1:A:143:C:C6	2.50	0.46

There are no symmetry-related clashes.

5.3 Torsion angles (i)

5.3.1 Protein backbone (i)

There are no protein molecules in this entry.

5.3.2 Protein sidechains (i)

There are no protein molecules in this entry.

5.3.3 RNA (i)

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
1	A	173/174 (99%)	18 (10%)	4 (2%)

5 of 18 RNA backbone outliers are listed below:



Mol	Chain	Res	Type
1	A	14	G
1	A	39	A
1	A	40	U
1	A	49	G
1	A	53	С

All (4) RNA pucker outliers are listed below:

Mol	Chain	Res	Type
1	A	39	A
1	A	52	U
1	A	53	С
1	A	80	G

5.4 Non-standard residues in protein, DNA, RNA chains (i)

1 non-standard protein/DNA/RNA residue is modelled in this entry.

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).

Mal	Tuno	Chain	Dec	Tiple	Bo	ond leng	$ ag{ths}$	B	ond ang	gles
MIOI	туре	Chain	nes	Lilik	Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z >2
1	CCC	A	174	1	20,25,26	1.01	1 (5%)	28,38,41	2.23	10 (35%)

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
1	CCC	A	174	1	-	2/7/35/36	0/3/3/3

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	\mathbf{Z}	$\operatorname{Observed}(\text{\AA})$	$Ideal(\AA)$
1	A	174	CCC	C6-C5	2.01	1.39	1.35



The worst	5	of	10	bond	angle	outliers	are	listed	below:
THE WOLDS	\circ	OI	10	DOM	angic	Outilities	COL C	mouca	DCIOW.

Mol	Chain	Res	Type	Atoms	\mathbf{Z}	$Observed(^o)$	$\operatorname{Ideal}({}^{o})$
1	A	174	CCC	O3'-PC-O1C	6.64	133.28	115.76
1	A	174	CCC	O2'-C2'-C1'	3.99	122.86	111.43
1	A	174	CCC	O3'-C3'-C2'	3.64	111.76	105.08
1	A	174	CCC	O2C-PC-O1C	3.50	121.18	109.89
1	A	174	CCC	O4'-C4'-C3'	-2.80	98.87	104.87

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	A	174	CCC	C3'-C4'-C5'-O5'
1	A	174	CCC	O4'-C4'-C5'-O5'

There are no ring outliers.

No monomer is involved in short contacts.

5.5 Carbohydrates (i)

There are no monosaccharides in this entry.

5.6 Ligand geometry (i)

Of 38 ligands modelled in this entry, 33 are monoatomic - leaving 5 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	Tuno	Chain	Res	Link	В	Bond lengths			Bond angles		
MIOI	Type	Chain	nes	Lilik	Counts	RMSZ	# Z >2	Counts	RMSZ	# Z > 2	
6	1PE	A	251	-	2,2,15	0.46	0	1,1,14	0.29	0	
6	1PE	A	252	-	4,4,15	0.56	0	3,3,14	0.29	0	
6	1PE	A	250	-	2,2,15	0.44	0	1,1,14	0.36	0	
7	IPA	A	253	-	3,3,3	0.54	0	3,3,3	0.20	0	
2	LYS	A	175	3	8,9,9	0.79	1 (12%)	9,10,10	1.44	2 (22%)	



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
6	1PE	A	252	-	-	2/2/2/13	_
2	LYS	A	175	3	-	0/9/9/9	-

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	\mathbf{Z}	$\operatorname{Observed}(\text{\AA})$	$Ideal(\AA)$
2	A	175	LYS	OXT-C	-2.03	1.23	1.30

All (2) bond angle outliers are listed below:

N	Mol	Chain	Res	Type	Atoms	Z	$Observed(^o)$	$\operatorname{Ideal}({}^{o})$
	2	A	175	LYS	OXT-C-O	-3.32	116.54	124.09
	2	A	175	LYS	OXT-C-CA	2.62	122.31	113.38

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
6	A	252	1PE	C13-C23-OH3-C22
6	A	252	1PE	C12-C22-OH3-C23

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mo	Chain	Res	Type	Clashes	Symm-Clashes
2	A	175	LYS	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, 95^{th} 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>	#RS	SRZ:	>2	$OWAB(Å^2)$	Q<0.9
1	A	173/174 (99%)	0.30	8 (4%)	32	35	13, 20, 42, 49	0

The worst 5 of 8 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	60	G	4.0
1	A	152	С	3.7
1	A	39	A	3.2
1	A	34	G	2.8
1	A	38	С	2.7

6.2 Non-standard residues in protein, DNA, RNA chains (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, 95^{th} 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	$\mathbf{B} ext{-}\mathbf{factors}(\mathbf{\mathring{A}}^2)$	Q < 0.9
1	CCC	A	174	23/24	0.83	0.23	26,29,31,31	0

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, 95^{th} 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	$\operatorname{B-factors}(\mathring{\mathbf{A}}^2)$	Q < 0.9
6	1PE	A	250	3/16	0.41	0.68	55,55,55,55	0
5	MG	A	229	1/1	0.69	0.31	39,39,39,39	0
7	IPA	A	253	4/4	0.69	0.29	42,42,42,42	0
4	NA	A	223	1/1	0.81	0.13	31,31,31,31	0
6	1PE	A	252	5/16	0.83	0.28	38,38,38,38	0
6	1PE	A	251	3/16	0.83	0.27	29,29,29,30	0
4	NA	A	226	1/1	0.86	0.14	35,35,35,35	0
4	NA	A	224	1/1	0.87	0.15	30,30,30,30	0
3	K	A	231	1/1	0.90	0.36	48,48,48,48	0
4	NA	A	222	1/1	0.90	0.14	25,25,25,25	0
4	NA	A	225	1/1	0.90	0.20	16,16,16,16	0
4	NA	A	215	1/1	0.91	0.13	37,37,37,37	0
4	NA	A	221	1/1	0.91	0.25	30,30,30,30	0
4	NA	A	214	1/1	0.93	0.09	35,35,35,35	0
4	NA	A	228	1/1	0.94	0.28	27,27,27,27	0
4	NA	A	217	1/1	0.94	0.24	30,30,30,30	0
4	NA	A	210	1/1	0.95	0.27	20,20,20,20	0
4	NA	A	218	1/1	0.95	0.17	25,25,25,25	0
4	NA	A	216	1/1	0.96	0.09	20,20,20,20	0
2	LYS	A	175	10/10	0.96	0.12	7,7,8,8	0
3	K	A	230	1/1	0.96	0.22	44,44,44,44	0
4	NA	A	220	1/1	0.96	0.10	23,23,23,23	0
4	NA	A	201	1/1	0.97	0.08	12,12,12,12	0
4	NA	A	211	1/1	0.98	0.05	16,16,16,16	0
3	K	A	200	1/1	0.98	0.09	13,13,13,13	0
4	NA	A	202	1/1	0.98	0.07	13,13,13,13	0
4	NA	A	203	1/1	0.98	0.08	19,19,19,19	0
4	NA	A	227	1/1	0.98	0.21	20,20,20,20	0
4	NA	A	204	1/1	0.98	0.10	15,15,15,15	0
4	NA	A	205	1/1	0.98	0.10	16,16,16,16	0
4	NA	A	219	1/1	0.98	0.11	31,31,31,31	0
4	NA	A	207	1/1	0.98	0.06	11,11,11,11	0
4	NA	A	209	1/1	0.98	0.15	11,11,11,11	0
4	NA	A	812	1/1	0.98	0.13	7,7,7,7	0
4	NA	A	208	1/1	0.99	0.05	19,19,19,19	0
4	NA	A	206	1/1	0.99	0.11	14,14,14,14	0
4	NA	A	212	1/1	0.99	0.23	15,15,15,15	0
4	NA	A	213	1/1	0.99	0.14	25,25,25,25	0



6.5 Other polymers (i)

There are no such residues in this entry.

