

wwPDB X-ray Structure Validation Summary Report (i)

May 21, 2020 – 12:59 am BST

PDB ID : 4ALP

Title: The Lin28b Cold shock domain in complex with hexauridine

Authors: Mayr, F.; Schuetz, A.; Doege, N.; Heinemann, U.

Deposited on : 2012-03-05

Resolution : 1.48 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 following versions of software and data (see references (1)) 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 : 2.11

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)

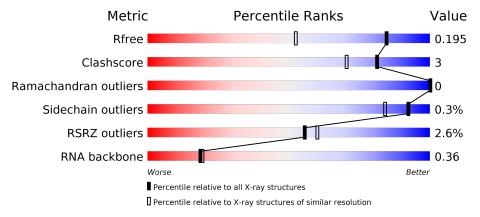
Validation Pipeline (wwPDB-VP) : 2.11

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 1.48 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	$\begin{array}{c} \text{Whole archive} \\ (\#\text{Entries}) \end{array}$	$\begin{array}{c} {\rm Similar \; resolution} \\ (\#{\rm Entries, \; resolution \; range(\AA)}) \end{array}$
R_{free}	130704	4690 (1.50-1.46)
Clashscore	141614	4955 (1.50-1.46)
Ramachandran outliers	138981	4846 (1.50-1.46)
Sidechain outliers	138945	4844 (1.50-1.46)
RSRZ outliers	127900	4614 (1.50-1.46)
RNA backbone	3102	1000 (2.34-0.62)

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 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					
1	A	90	96					
1	Λ	90	94% 3%	• •				
1	В	90	91%					
1	C	00	2%					
1	C	90	91%	• •				
1	D	90	92%	• 7%				

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Mol	Chain	Length	Quality of chain						
2	Е	6	33%	33%	17%	17%			



2 Entry composition (i)

There are 4 unique types of molecules in this entry. The entry contains 3566 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 LIN28 ISOFORM B.

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace		
1	Λ	0.7	Total	С	N	О	S	0	4	0
1	A	87	691	438	120	128	5	0	$\frac{4}{2}$	0
1	В	86	Total	С	N	О	S	0	4	0
1	Ъ	00	692	439	121	127	5		'±	. 0
1	С	86	Total	С	N	О	S	0	7	0
1		00	713	454	124	129	6			
1	D	84	Total	С	N	О	S	0	5	0
1			687	437	124	121	5			

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference	
A	25	GLY	-	expression tag	UNP B4F6I0	
A	26	SER	_	expression tag	UNP B4F6I0	
В	25	GLY	-	expression tag	UNP B4F6I0	
В	26	SER	=	expression tag	UNP B4F6I0	
С	25	GLY	-	expression tag	UNP B4F6I0	
С	26	SER	-	expression tag	UNP B4F6I0	
D	25	GLY	=	expression tag	UNP B4F6I0	
D	26	SER	-	expression tag	UNP B4F6I0	

• Molecule 2 is a RNA chain called HEXA URIDINE.

Mol	Chain	Residues	${f Atoms}$			ZeroOcc	AltConf	Trace		
2	Е	6	Total 115	C 53	N 12	O 45	P 5	0	0	0

• Molecule 3 is GLYCEROL (three-letter code: GOL) (formula: C₃H₈O₃).





Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	В	1	Total C O 6 3 3	0	0
3	D	1	Total C O 6 3 3	0	0

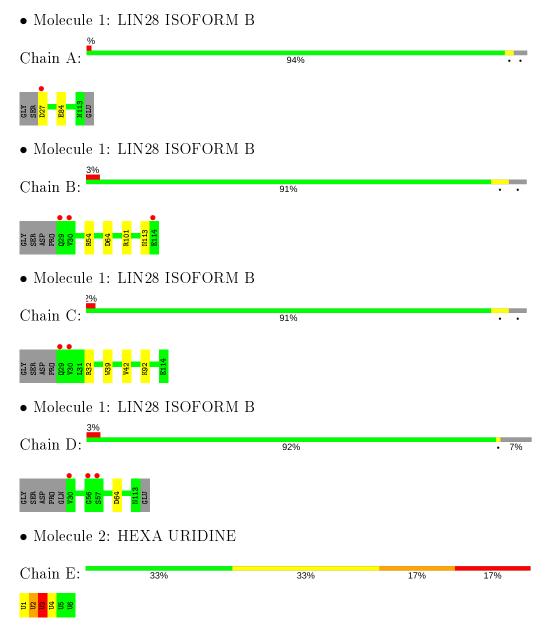
• Molecule 4 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	201	Total O 201 201	0	0
4	В	149	Total O 149 149	0	0
4	С	153	Total O 153 153	0	0
4	D	127	Total O 127 127	0	0
4	Е	26	Total O 26 26	0	0



3 Residue-property plots (i)

These plots are drawn for all protein, RNA and DNA 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.





4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants	95.98Å 61.17Å 63.22Å	Danasitan
a, b, c, α , β , γ	90.00° 90.00° 90.00°	Depositor
Resolution (Å)	32.42 - 1.48	Depositor
Resolution (A)	29.14 - 1.48	EDS
% Data completeness	96.4 (32.42-1.48)	Depositor
(in resolution range)	96.4 (29.14-1.48)	EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) > 1$	3.50 (at 1.48Å)	Xtriage
Refinement program	REFMAC 5.6.0117	Depositor
D D	0.169 , 0.188	Depositor
R, R_{free}	0.178 , 0.195	DCC
R_{free} test set	3023 reflections $(5.00%)$	wwPDB-VP
Wilson B-factor (Å ²)	11.6	Xtriage
Anisotropy	0.046	Xtriage
Bulk solvent $k_{sol}(e/Å^3)$, $B_{sol}(Å^2)$	0.32 , 40.4	EDS
L-test for twinning ²	$< L >=0.47, < L^2>=0.30$	Xtriage
Estimated twinning fraction	0.025 for -h,l,k	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	3566	wwPDB-VP
Average B, all atoms (Å ²)	16.0	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 21.47 % 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 6.9659e-03.

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: GOL

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	Bond angles		
MIOI	Chain	RMSZ	# Z > 5	RMSZ	# Z > 5	
1	A	0.74	0/718	0.75	0/963	
1	В	0.68	0/721	0.81	2/964 (0.2%)	
1	С	0.72	1/744~(0.1%)	0.79	0/992	
1	D	0.69	0/719	0.79	1/960 (0.1%)	
2	Е	0.69	0/125	1.46	4/190 (2.1%)	
All	All	0.71	1/3027 (0.0%)	0.83	7/4069 (0.2%)	

All (1) bond length outliers are listed below:

\mathbf{Mol}	Chain	Res	Type	Atoms	Z	$\operatorname{Observed}(\operatorname{\AA})$	$\operatorname{Ideal}(ext{\AA})$
1	С	39	TRP	CD2-CE2	5.74	1.48	1.41

The worst 5 of 7 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	\mathbf{Z}	$\mathbf{Observed}(^o)$	$\mathbf{Ideal}(^o)$
2	Е	2	U	O3'-P-O5'	-9.18	86.56	104.00
2	E	3	U	O5'-P-OP2	7.31	119.47	110.70
1	В	64	ASP	CB-CG-OD1	6.35	124.02	118.30
2	Е	3	U	P-O3'-C3'	6.22	127.17	119.70
1	В	64	ASP	CB-CG-OD2	-6.03	112.87	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
--

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	691	0	675	1	0
1	В	692	0	682	5	0
1	С	713	0	712	3	0
1	D	687	0	689	0	0
2	E	115	0	61	3	0
3	В	6	0	8	3	0
3	D	6	0	8	0	0
4	A	201	0	0	1	0
4	В	149	0	0	5	0
4	С	153	0	0	2	0
4	D	127	0	0	0	0
4	E	26	0	0	1	0
All	All	3566	0	2835	15	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 15 close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	$egin{aligned} ext{Interatomic} \ ext{distance} \ (ext{Å}) \end{aligned}$	$egin{array}{c} ext{Clash} \ ext{overlap } (ext{Å}) \end{array}$
1:B:54[B]:ARG:HG2	1:B:54[B]:ARG:HH11	1.31	0.94
1:B:101:ARG:HD2	4:B:2130:HOH:O	1.73	0.88
1:B:54[B]:ARG:HH11	1:B:54[B]:ARG:CG	1.87	0.86
3:B:1115:GOL:C1	4:B:2027:HOH:O	2.30	0.78
2:E:3:U:N3	2:E:3:U:N1	2.28	0.73

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	89/90 (99%)	86 (97%)	3 (3%)	0	100	100
1	В	88/90 (98%)	86 (98%)	2 (2%)	0	100	100
1	С	90/90~(100%)	89 (99%)	1 (1%)	0	100	100
1	D	87/90 (97%)	86 (99%)	1 (1%)	0	100	100
All	All	$354/360 \ (98\%)$	347 (98%)	7 (2%)	0	100	100

There are no Ramachandran outliers to report.

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	78/77 (101%)	77 (99%)	1 (1%)	69	42	
1	В	78/77~(101%)	78 (100%)	0	100	100	
1	С	81/77 (105%)	81 (100%)	0	100	100	
1	D	77/77~(100%)	77 (100%)	0	100	100	
All	All	314/308 (102%)	313 (100%)	1 (0%)	92	84	

All (1) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	84	GLU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	С	113	ASN

5.3.3 RNA (i)



Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
2	E	5/6 (83%)	3 (60%)	1 (20%)

All (3) RNA backbone outliers are listed below:

Mol	Chain	Res	Type
2	Е	2	U
2	E	3	U
2	Е	4	U

All (1) RNA pucker outliers are listed below:

Mol	Chain	Res	Type
2	Е	3	U

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 carbohydrates in this entry.

5.6 Ligand geometry (i)

2 ligands are 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).

Mol	Tree o	Chain	Dag	Timl.	В	ond leng	$_{ m gths}$	В	ond ang	gles
MIOI	Type	Chain	Res	Link	Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	GOL	D	1114	-	5,5,5	0.29	0	5, 5, 5	0.36	0
3	GOL	В	1115	-	5,5,5	0.82	0	5,5,5	1.01	0

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
3	GOL	D	1114	_	-	0/4/4/4	ı
3	GOL	В	1115	-	-	3/4/4/4	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (3) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	В	1115	GOL	O1-C1-C2-O2
3	В	1115	GOL	O1-C1-C2-C3
3	В	1115	GOL	C1-C2-C3-O3

There are no ring outliers.

1 monomer is involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	В	1115	GOL	3	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	$\langle { m RSRZ} \rangle$	$\# \mathrm{RSRZ}{>}2$	$OWAB(\AA^2)$	Q < 0.9
1	A	87/90 (96%)	-0.09	1 (1%) 80 83	7, 10, 21, 38	0
1	В	86/90 (95%)	0.04	3 (3%) 44 48	7, 11, 29, 41	0
1	С	86/90 (95%)	-0.03	2 (2%) 60 65	8, 12, 22, 39	0
1	D	84/90 (93%)	0.09	3 (3%) 42 46	7, 16, 29, 40	0
2	E	6/6 (100%)	0.35	0 100 100	10, 15, 23, 26	0
All	All	349/366 (95%)	0.01	9 (2%) 56 60	7, 12, 27, 41	0

The worst 5 of 9 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	С	29	GLN	4.4
1	С	30	VAL	4.1
1	В	29	GLN	3.3
1	В	30	VAL	3.2
1	D	56	GLY	3.0

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. 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
3	GOL	В	1115	6/6	0.83	0.21	14,21,31,36	0
3	GOL	D	1114	6/6	0.92	0.21	27,28,29,32	0

6.5 Other polymers (i)

There are no such residues in this entry.

