

wwPDB X-ray Structure Validation Summary Report (i)

Aug 20, 2023 - 08:30 AM EDT

PDB ID : 2LYN

Title : HIGH RESOLUTION STRUCTURE OF RED ABALONE LYSIN DIMER

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Deposited on : 1999-04-19

Resolution : 2.07 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:

 $\begin{array}{ccc} \text{MolProbity} & : & 4.02\text{b-}467 \\ \text{Xtriage (Phenix)} & : & 1.13 \end{array}$

(Phenix) : 1.13 EDS : 2.35

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)

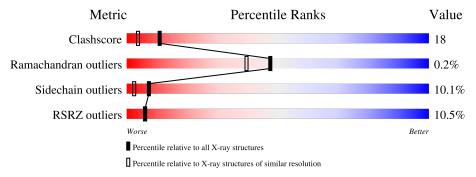
Validation Pipeline (wwPDB-VP) : 2.35

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 2.07 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	Similar resolution
Metric	$(\# \mathrm{Entries})$	$(\# ext{Entries}, ext{ resolution range}(ext{Å}))$
Clashscore	141614	2801 (2.08-2.04)
Ramachandran outliers	138981	2768 (2.08-2.04)
Sidechain outliers	138945	2768 (2.08-2.04)
RSRZ outliers	127900	2646 (2.08-2.04)

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				
1	A	136	6%	8%	23%	5% • •		
1	В	136	28%	46%		10% 7%		
1	С	136	2%	74%	16%	• 6%		
1	D	136	4%	74%	15%	• • 5%		



2 Entry composition (i)

There are 2 unique types of molecules in this entry. The entry contains 4618 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 SPERM LYSIN.

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace			
1	٨	131	Total	С	N	О	S	0	1	1 1	
1	A	131	1101	720	195	179	7	0	1	1	
1	В	126	Total	С	N	О	S	0	0	1	
1	Б	120	1051	687	186	171	7	0	U	1	
1	С	128	Total	С	N	О	S	0	1	1	
1		120	1074	702	192	173	7	0	1	1	
1	D	129	Total	С	N	О	S	0	0	1	
1	D	129	1076	703	190	176	7	U	U	1	

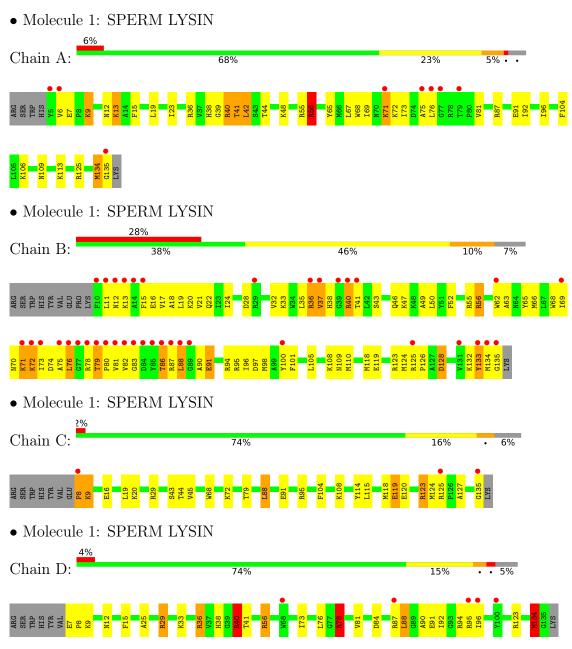
• Molecule 2 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	93	Total O 93 93	0	0
2	В	38	Total O 38 38	0	0
2	С	93	Total O 93 93	0	0
2	D	92	Total O 92 92	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.





4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 41 21 2	Depositor
Cell constants	72.52Å 72.52Å 266.08Å	Donositor
a, b, c, α , β , γ	90.00° 90.00° 90.00°	Depositor
Resolution (Å)	100.00 - 2.07	Depositor
resolution (A)	37.83 - 2.07	EDS
% Data completeness	99.1 (100.00-2.07)	Depositor
(in resolution range)	99.4 (37.83-2.07)	EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.09	Depositor
$< I/\sigma(I) > 1$	3.14 (at 2.08Å)	Xtriage
Refinement program	SHELXL	Depositor
P. P.	0.216 , 0.285	Depositor
R, R_{free}	0.205 , (Not available)	DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å ²)	26.3	Xtriage
Anisotropy	0.006	Xtriage
Bulk solvent $k_{sol}(e/Å^3)$, $B_{sol}(Å^2)$	0.36 , 111.8	EDS
L-test for twinning ²	$ < L >=0.47, < 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	4618	wwPDB-VP
Average B, all atoms (Å ²)	37.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: The largest off-origin peak in the Patterson function is 8.10% 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)

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).

Mal	Mol Chain		nd lengths	Bond angles		
IVIOI			# Z > 5	RMSZ	# Z > 5	
1	A	0.56	1/1135 (0.1%)	1.07	$6/1530 \ (0.4\%)$	
1	В	0.26	0/1077	0.80	0/1451	
1	С	0.30	0/1105	0.88	$2/1487 \ (0.1\%)$	
1	D	0.59	1/1104 (0.1%)	0.98	8/1488 (0.5%)	
All	All	0.45	$2/4421 \ (0.0\%)$	0.94	$16/5956 \ (0.3\%)$	

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	\mathbf{Z}	Observed(A)	$\operatorname{Ideal}(ext{\AA})$
1	D	134	MET	C-N	-16.46	1.03	1.33
1	A	134	MET	C-N	15.86	1.61	1.33

The worst 5 of 16 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	\mathbf{Z}	$Observed(^o)$	$Ideal(^{o})$
1	A	56	ARG	CD-NE-CZ	22.71	155.40	123.60
1	A	56	ARG	NE-CZ-NH2	8.40	124.50	120.30
1	С	8	PRO	C-N-CA	6.72	138.50	121.70
1	D	40	ARG	NE-CZ-NH1	-6.70	116.95	120.30
1	A	134	MET	O-C-N	-6.67	111.86	123.20

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	1101	0	1132	29	0
1	В	1051	0	1079	76	0
1	С	1074	0	1109	16	0
1	D	1076	0	1105	40	0
2	A	93	0	0	2	0
2	В	38	0	0	4	0
2	С	93	0	0	3	0
2	D	92	0	0	3	0
All	All	4618	0	4425	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 18.

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	Clash
710011-1	1100111-2	${ m distance}({ m \AA})$	overlap (Å)
1:D:91:GLU:HG2	1:D:94:ARG:HH21	1.16	1.08
1:D:91:GLU:HA	1:D:94:ARG:HE	1.44	0.81
1:A:44:THR:HG22	1:A:48:LYS:HZ2	1.46	0.80
1:A:75:ALA:HB2	2:A:220:HOH:O	1.83	0.78
1:D:7:GLU:OE2	1:D:8:PRO:HD2	1.84	0.78

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	Perce	ntiles
1	A	130/136 (96%)	128 (98%)	2 (2%)	0	100	100
1	В	123/136 (90%)	115 (94%)	7 (6%)	1 (1%)	19	9
1	С	126/136 (93%)	124 (98%)	2 (2%)	0	100	100
1	D	127/136 (93%)	125 (98%)	2 (2%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
All	All	506/544 (93%)	492 (97%)	13 (3%)	1 (0%)	47 39

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	В	76	LEU

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	Perc	entiles
1	A	115/119 (97%)	101 (88%)	14 (12%)	5	1
1	В	109/119 (92%)	93 (85%)	16 (15%)	3	0
1	С	112/119 (94%)	105 (94%)	7 (6%)	18	10
1	D	112/119 (94%)	104 (93%)	8 (7%)	14	7
All	All	448/476 (94%)	403 (90%)	45 (10%)	7	2

5 of 45 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	В	128	ASP
1	С	119	GLU
1	В	133	TYR
1	С	44	THR
1	D	33	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 6 such sidechains are listed below:

Mol	Chain	Res	Type
1	С	64	ASN
1	С	70	ASN
1	D	70	ASN
1	В	70	ASN

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Mol	Chain	Res	Type
1	В	64	ASN

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)

The following chains have linkage breaks:

Mol	Chain	Number of breaks
1	В	1
1	С	1
1	A	1
1	D	1

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	В	134:MET	С	135:GLY	N	5.11
1	С	134:MET	С	135:GLY	N	3.33
1	A	134:MET	С	135:GLY	N	1.61

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Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	D	134:MET	С	135:GLY	N	1.03



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 > 2	$OWAB(Å^2)$	Q<0.9
1	A	131/136 (96%)	0.37	8 (6%) 21 21	17, 28, 64, 89	0
1	В	126/136 (92%)	1.60	38 (30%) 0 0	25, 51, 111, 131	0
1	С	128/136 (94%)	0.02	3 (2%) 60 63	17, 28, 52, 99	0
1	D	129/136 (94%)	0.29	5 (3%) 39 41	15, 25, 61, 82	0
All	All	514/544 (94%)	0.56	54 (10%) 6 6	15, 30, 79, 131	0

The worst 5 of 54 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	В	76	LEU	8.4
1	В	81	VAL	7.0
1	В	79	THR	6.7
1	В	73	ILE	6.2
1	A	135	GLY	6.0

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 monosaccharides in this entry.

6.4 Ligands (i)

There are no ligands in this entry.



6.5 Other polymers (i)

There are no such residues in this entry.

