

wwPDB X-ray Structure Validation Summary Report (i)

Dec 16, 2023 – 03:12 PM EST

PDB ID : 2OQA

Title : X-ray Sequence and Crystal Structure of Luffaculin 1, a Novel Type 1 Riboso

me-inactivating Protein

Authors : Hou, X.; Huang, M.

Deposited on : 2007-01-31

Resolution : 1.40 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)

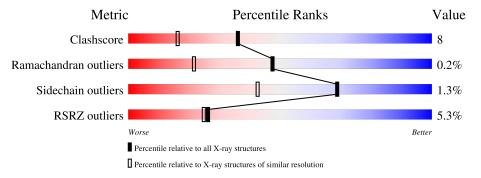
Validation Pipeline (wwPDB-VP) : 2.36

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 1.40 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.



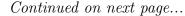
Metric	Whole archive	Similar resolution
	$(\# \mathrm{Entries})$	$(\# ext{Entries}, ext{ resolution range}(ext{Å}))$
Clashscore	141614	1812 (1.40-1.40)
Ramachandran outliers	138981	1763 (1.40-1.40)
Sidechain outliers	138945	1762 (1.40-1.40)
RSRZ outliers	127900	1674 (1.40-1.40)

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	241	87%	12%			
1	В	241	7% 88%	7% • •			

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
2	NAG	A	1001	-	-	X	-
2	NAG	A	1004	-	-	-	X





Continued from previous page...

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	PG4	A	801	-	-	-	X



2 Entry composition (i)

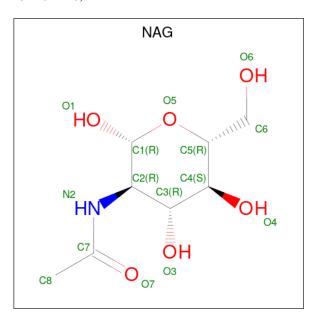
There are 5 unique types of molecules in this entry. The entry contains 4177 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Luffaculin 1.

\mathbf{Mol}	Chain	Residues	${f Atoms}$			ZeroOcc	AltConf	Trace		
1	A	241	Total		- '	0	S	0	0	0
			1825	1165	299	357	4			
1	B	233	Total	\mathbf{C}	N	O	\mathbf{S}	0	0	0
1	Ъ	255	1775	1138	289	345	3		0	0

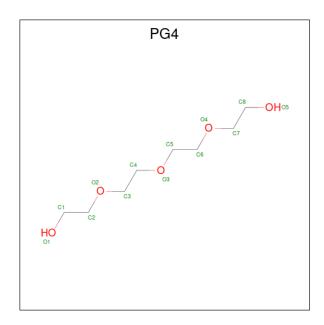
• Molecule 2 is 2-acetamido-2-deoxy-beta-D-glucopyranose (three-letter code: NAG) (formula: $C_8H_{15}NO_6$).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total C N O 15 8 1 6	0	0
2	A	1	Total C N O 15 8 1 6	0	0
2	В	1	Total C N O 15 8 1 6	0	0
2	В	1	Total C N O 15 8 1 6	0	0

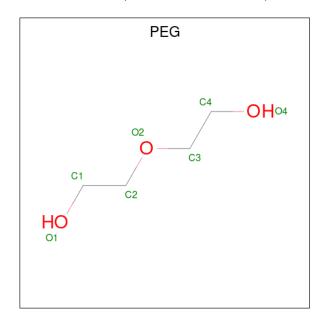


• Molecule 3 is TETRAETHYLENE GLYCOL (three-letter code: PG4) (formula: $C_8H_{18}O_5$).



Mol	Chain	Residues	Aton	ns	ZeroOcc	AltConf
3	A	1	Total (C O 8 5	0	0

• Molecule 4 is DI(HYDROXYETHYL)ETHER (three-letter code: PEG) (formula: $C_4H_{10}O_3$).



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



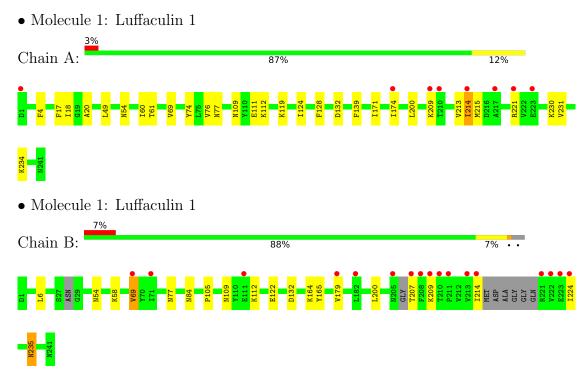
• Molecule 5 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	248	Total O 248 248	0	0
5	В	242	Total O 242 242	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 1	Depositor
Cell constants	39.13Å 46.81Å 83.57Å	Depositor
a, b, c, α , β , γ	89.07° 80.01° 72.14°	Depositor
Resolution (Å)	33.90 - 1.40	Depositor
resolution (A)	33.90 - 1.40	EDS
% Data completeness	86.7 (33.90-1.40)	Depositor
(in resolution range)	86.6 (33.90-1.40)	EDS
R_{merge}	0.03	Depositor
R_{sym}	0.12	Depositor
$< I/\sigma(I) > 1$	5.41 (at 1.40Å)	Xtriage
Refinement program	REFMAC 5.2.0005	Depositor
R, R_{free}	0.213 , 0.232	Depositor
it, it free	0.215 , (Not available)	DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor (A^2)	13.1	Xtriage
Anisotropy	0.688	Xtriage
Bulk solvent $k_{sol}(e/Å^3)$, $B_{sol}(Å^2)$	0.40, 46.8	EDS
L-test for twinning ²	$ < L > = 0.49, < L^2> = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	4177	wwPDB-VP
Average B, all atoms $(Å^2)$	18.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: The largest off-origin peak in the Patterson function is 6.88% 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: NAG, PG4, PEG

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		
MIOI	Chain	RMSZ	# Z > 5	RMSZ	# Z > 5	
1	A	0.37	0/1855	0.53	0/2515	
1	В	0.38	0/1802	0.52	0/2440	
All	All	0.38	0/3657	0.53	0/4955	

There are no bond length outliers.

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	1825	0	1860	37	0
1	В	1775	0	1816	22	0
2	A	30	0	30	9	0
2	В	30	0	30	10	0
3	A	13	0	18	0	0
4	В	14	0	20	0	0
5	A	248	0	0	12	0
5	В	242	0	0	4	0
All	All	4177	0	3774	59	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 8.



The worst 5 of 59 close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	$\begin{array}{c} \text{Interatomic} \\ \text{distance (Å)} \end{array}$	$\begin{array}{c} \text{Clash} \\ \text{overlap } (\text{\AA}) \end{array}$
1:B:77:ASN:ND2	2:B:1002:NAG:H1	1.28	1.46
1:A:77:ASN:HD21	2:A:1001:NAG:C1	1.45	1.30
1:B:77:ASN:HD21	2:B:1002:NAG:C1	1.58	1.16
1:B:77:ASN:ND2	2:B:1002:NAG:C1	2.09	1.14
1:A:77:ASN:ND2	2:A:1001:NAG:C1	2.26	0.96

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	$_{ m ntiles}$
1	A	239/241 (99%)	229 (96%)	10 (4%)	0	100	100
1	В	225/241 (93%)	216 (96%)	8 (4%)	1 (0%)	34	12
All	All	464/482 (96%)	445 (96%)	18 (4%)	1 (0%)	47	21

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	В	69	VAL

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	Percentile	s
1	A	201/201 (100%)	199 (99%)	2 (1%)	76 53	
1	В	197/201 (98%)	194 (98%)	3 (2%)	65 37	
All	All	398/402 (99%)	393 (99%)	5 (1%)	69 42	

All (5) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	69	VAL
1	A	214	ILE
1	В	69	VAL
1	В	164	LYS
1	В	235	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 10 such sidechains are listed below:

Mol	Chain	Res	Type
1	В	205	ASN
1	В	235	ASN
1	В	237	GLN
1	A	220	GLN
1	В	77	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)

7 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	Trino	Chain	Res	Link	Вс	ond leng	ths	В	ond ang	les
MIOI	Type	Chain	nes	LIIIK	Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
4	PEG	В	802	-	6,6,6	0.43	0	5,5,5	0.30	0
4	PEG	В	803	-	6,6,6	0.44	0	5,5,5	0.32	0
2	NAG	В	1002	-	15,15,15	0.80	0	21,21,21	2.96	4 (19%)
2	NAG	A	1001	-	15,15,15	0.73	0	21,21,21	3.78	7 (33%)
3	PG4	A	801	-	12,12,12	0.54	0	11,11,11	0.30	0
2	NAG	A	1004	-	15,15,15	0.42	0	21,21,21	0.67	0
2	NAG	В	1003	1	15,15,15	0.54	0	21,21,21	0.92	1 (4%)

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
4	PEG	В	802	-	-	0/4/4/4	-
4	PEG	В	803	-	-	3/4/4/4	-
2	NAG	В	1002	-	-	2/6/26/26	0/1/1/1
2	NAG	A	1001	-	-	1/6/26/26	0/1/1/1
3	PG4	A	801	-	-	3/10/10/10	-
2	NAG	A	1004	-	-	4/6/26/26	0/1/1/1
2	NAG	В	1003	1	-	3/6/26/26	0/1/1/1

There are no bond length outliers.

The worst 5 of 12 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	\mathbf{Z}	$Observed(^o)$	$\operatorname{Ideal}({}^{o})$
2	A	1001	NAG	C1-C2-C3	-13.56	92.05	110.54
2	В	1002	NAG	C1-C2-N2	9.91	122.22	110.73
2	A	1001	NAG	C1-O5-C5	-7.95	98.67	113.66
2	В	1002	NAG	O5-C1-C2	7.49	117.04	109.52
2	A	1001	NAG	O1-C1-O5	3.84	121.91	110.38



There are no chirality outliers.

5 of 16 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	1004	NAG	C8-C7-N2-C2
2	A	1004	NAG	O7-C7-N2-C2
2	В	1003	NAG	C8-C7-N2-C2
2	В	1003	NAG	O7-C7-N2-C2
2	A	1004	NAG	O5-C5-C6-O6

There are no ring outliers.

3 monomers are involved in 19 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	В	1002	NAG	6	0
2	A	1001	NAG	9	0
2	В	1003	NAG	4	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 $>$	$\# \mathrm{RSRZ}{>}2$	$OWAB(Å^2)$	Q<0.9
1	A	241/241 (100%)	0.28	8 (3%) 46 46	12, 16, 25, 30	0
1	В	$233/241 \ (96\%)$	0.51	17 (7%) 15 13	10, 16, 27, 40	0
All	All	474/482 (98%)	0.39	25 (5%) 26 25	10, 16, 26, 40	0

The worst 5 of 25 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	В	214	ILE	4.2
1	В	182	LEU	4.0
1	В	221	ARG	3.5
1	В	205	ASN	2.9
1	В	209	LYS	2.9

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, 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
2	NAG	A	1004	15/15	0.33	0.51	40,45,46,46	0
3	PG4	A	801	13/13	0.40	0.41	69,69,69,69	0
4	PEG	В	803	7/7	0.59	0.32	52,53,53,53	0
4	PEG	В	802	7/7	0.78	0.26	38,39,39,40	0
2	NAG	В	1003	15/15	0.83	0.17	25,26,28,30	0
2	NAG	В	1002	15/15	0.85	0.15	13,14,19,25	0
2	NAG	A	1001	15/15	0.88	0.14	12,15,18,25	0

6.5 Other polymers (i)

There are no such residues in this entry.

