



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 13, 2023 – 10:07 AM EDT

PDB ID : 8FWL
Title : Crystal structure of Australian bat lyssavirus nucleoprotein in complex with phosphoprotein chaperone
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Deposited on : 2023-01-23
Resolution : 2.19 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 types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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 : 2.32.1
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) : 2.32.1

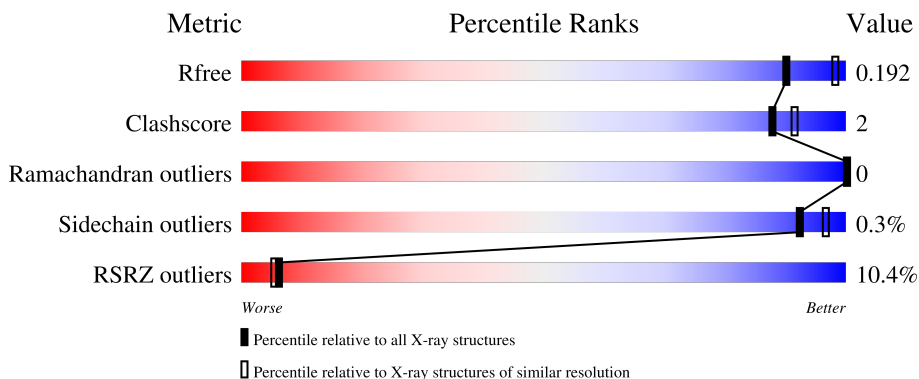
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.19 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	4898 (2.20-2.20)
Clashscore	141614	5594 (2.20-2.20)
Ramachandran outliers	138981	5503 (2.20-2.20)
Sidechain outliers	138945	5504 (2.20-2.20)
RSRZ outliers	127900	4800 (2.20-2.20)

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	567	

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 6747 atoms, of which 3264 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 Phosphoprotein,Nucleoprotein.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
			Total	C	H	N	O	S			
1	A	413	6562	2099	3254	570	618	21	0	6	0

There are 64 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	MET	-	expression tag	UNP A0A7M4C440
A	2	HIS	-	expression tag	UNP A0A7M4C440
A	3	HIS	-	expression tag	UNP A0A7M4C440
A	4	HIS	-	expression tag	UNP A0A7M4C440
A	5	HIS	-	expression tag	UNP A0A7M4C440
A	6	HIS	-	expression tag	UNP A0A7M4C440
A	7	HIS	-	expression tag	UNP A0A7M4C440
A	8	SER	-	expression tag	UNP A0A7M4C440
A	9	SER	-	expression tag	UNP A0A7M4C440
A	10	GLY	-	expression tag	UNP A0A7M4C440
A	11	LEU	-	expression tag	UNP A0A7M4C440
A	12	VAL	-	expression tag	UNP A0A7M4C440
A	13	PRO	-	expression tag	UNP A0A7M4C440
A	14	ARG	-	expression tag	UNP A0A7M4C440
A	15	GLY	-	expression tag	UNP A0A7M4C440
A	16	SER	-	expression tag	UNP A0A7M4C440
A	17	GLY	-	expression tag	UNP A0A7M4C440
A	18	MET	-	expression tag	UNP A0A7M4C440
A	19	LYS	-	expression tag	UNP A0A7M4C440
A	20	GLU	-	expression tag	UNP A0A7M4C440
A	21	THR	-	expression tag	UNP A0A7M4C440
A	22	ALA	-	expression tag	UNP A0A7M4C440
A	23	ALA	-	expression tag	UNP A0A7M4C440
A	24	ALA	-	expression tag	UNP A0A7M4C440
A	25	LYS	-	expression tag	UNP A0A7M4C440
A	26	PHE	-	expression tag	UNP A0A7M4C440
A	27	GLU	-	expression tag	UNP A0A7M4C440

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Chain	Residue	Modelled	Actual	Comment	Reference
A	28	ARG	-	expression tag	UNP A0A7M4C440
A	29	GLN	-	expression tag	UNP A0A7M4C440
A	30	HIS	-	expression tag	UNP A0A7M4C440
A	31	MET	-	expression tag	UNP A0A7M4C440
A	32	ASP	-	expression tag	UNP A0A7M4C440
A	33	SER	-	expression tag	UNP A0A7M4C440
A	34	PRO	-	expression tag	UNP A0A7M4C440
A	35	ASP	-	expression tag	UNP A0A7M4C440
A	36	LEU	-	expression tag	UNP A0A7M4C440
A	37	GLY	-	expression tag	UNP A0A7M4C440
A	38	THR	-	expression tag	UNP A0A7M4C440
A	39	ASP	-	expression tag	UNP A0A7M4C440
A	40	ASP	-	expression tag	UNP A0A7M4C440
A	41	ASP	-	expression tag	UNP A0A7M4C440
A	42	ASP	-	expression tag	UNP A0A7M4C440
A	43	LYS	-	expression tag	UNP A0A7M4C440
A	44	ALA	-	expression tag	UNP A0A7M4C440
A	45	MET	-	expression tag	UNP A0A7M4C440
A	46	ALA	-	expression tag	UNP A0A7M4C440
A	47	ASP	-	expression tag	UNP A0A7M4C440
A	48	ILE	-	expression tag	UNP A0A7M4C440
A	49	GLY	-	expression tag	UNP A0A7M4C440
A	50	SER	-	expression tag	UNP A0A7M4C440
A	51	GLU	-	expression tag	UNP A0A7M4C440
A	52	ASN	-	expression tag	UNP A0A7M4C440
A	53	LEU	-	expression tag	UNP A0A7M4C440
A	54	TYR	-	expression tag	UNP A0A7M4C440
A	55	PHE	-	expression tag	UNP A0A7M4C440
A	56	GLN	-	expression tag	UNP A0A7M4C440
A	57	SER	-	expression tag	UNP A0A7M4C440
A	138	GLU	-	linker	UNP A0A7M4C440
A	139	ASN	-	linker	UNP A0A7M4C440
A	140	LEU	-	linker	UNP A0A7M4C440
A	141	TYR	-	linker	UNP A0A7M4C440
A	142	PHE	-	linker	UNP A0A7M4C440
A	143	GLN	-	linker	UNP A0A7M4C440
A	144	GLY	-	linker	UNP A0A7M4C440

- Molecule 2 is DI(HYDROXYETHYL)ETHER (three-letter code: PEG) (formula: C₄H₁₀O₃).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	C	H	O		
2	A	1	17	4	10	3	0	0

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
3	A	168	168	168	0	0

4 Data and refinement statistics i

Property	Value	Source
Space group	P 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	82.37Å 35.50Å 89.09Å 90.00° 92.63° 90.00°	Depositor
Resolution (Å)	29.67 – 2.19 29.67 – 2.19	Depositor EDS
% Data completeness (in resolution range)	99.6 (29.67-2.19) 99.6 (29.67-2.19)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.59 (at 2.20Å)	Xtrriage
Refinement program	PHENIX (1.20.1_4487: ???)	Depositor
R, R_{free}	0.159 , 0.193 0.158 , 0.192	Depositor DCC
R_{free} test set	1371 reflections (5.10%)	wwPDB-VP
Wilson B-factor (Å ²)	30.9	Xtrriage
Anisotropy	0.420	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.40 , 56.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtrriage
Estimated twinning fraction	0.028 for h,-k,-l	Xtrriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	6747	wwPDB-VP
Average B, all atoms (Å ²)	49.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.64% 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: 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	
		RMSZ	# Z >5	RMSZ	# Z >5
1	A	0.28	0/3382	0.49	0/4571

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	3308	3254	3245	13	0
2	A	7	10	10	0	0
3	A	168	0	0	1	0
All	All	3483	3264	3255	13	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 2.

All (13) 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:A:231:GLU:OE1	3:A:701:HOH:O	2.14	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:234:THR:OG1	1:A:239[A]:MET:SD	2.64	0.56
1:A:239[A]:MET:HE3	1:A:242:ARG:HG3	1.91	0.52
1:A:183:PRO:HG3	1:A:438:LEU:HD11	2.00	0.44
1:A:455:LEU:HD13	1:A:455:LEU:C	2.39	0.43
1:A:548:ALA:O	1:A:552:ARG:HG3	2.19	0.43
1:A:273:THR:HG22	1:A:274:ARG:N	2.34	0.43
1:A:405:PHE:HZ	1:A:553[B]:ILE:HD12	1.85	0.41
1:A:228:VAL:HG22	1:A:246:LYS:HD2	2.02	0.41
1:A:565:ILE:HG22	1:A:569:ILE:HD12	2.03	0.41
1:A:226:GLU:HG2	1:A:246:LYS:HE3	2.04	0.40
1:A:90[A]:GLU:OE2	1:A:391:LYS:NZ	2.51	0.40
1:A:575:HIS:CG	1:A:576:GLN:N	2.90	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	413/567 (73%)	407 (98%)	6 (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	358/487 (74%)	357 (100%)	1 (0%)	92 97

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

Mol	Chain	Res	Type
1	A	568	TYR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

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](#)

1 ligand 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).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
2	PEG	A	601	-	6,6,6	0.16	0	5,5,5	0.18	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
2	PEG	A	601	-	-	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
2	A	601	PEG	O2-C3-C4-O4
2	A	601	PEG	O1-C1-C2-O2
2	A	601	PEG	C1-C2-O2-C3

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	413/567 (72%)	0.57	43 (10%) 6 5	17, 39, 89, 130	0

All (43) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	301	ASN	10.1
1	A	300	GLN	8.4
1	A	268	GLY	8.4
1	A	272	LEU	7.2
1	A	60	LYS	7.0
1	A	544	ARG	6.8
1	A	267	THR	6.4
1	A	269	GLY	6.0
1	A	271	ASP	5.3
1	A	266	LEU	5.3
1	A	70	ALA	5.1
1	A	578	ARG	4.9
1	A	97	GLN	4.8
1	A	59	SER	4.5
1	A	74	ASP	4.5
1	A	496	LYS	4.4
1	A	576	GLN	4.4
1	A	577	SER	4.3
1	A	270	GLN	4.1
1	A	559	ARG	3.9
1	A	299	GLY	3.9
1	A	562	ARG	3.8
1	A	273	THR	3.8
1	A	574	ASN	3.5
1	A	61	ILE	3.3
1	A	68	ILE	3.2
1	A	72	MET	3.2

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Mol	Chain	Res	Type	RSRZ
1	A	495	GLY	2.9
1	A	274	ARG	2.9
1	A	303	GLY	2.9
1	A	557	GLY	2.8
1	A	69	ARG	2.7
1	A	67	ALA	2.7
1	A	591	SER	2.5
1	A	171	GLU	2.4
1	A	302[A]	THR	2.3
1	A	579	PRO	2.3
1	A	90[A]	GLU	2.2
1	A	73	ALA	2.2
1	A	289	LEU	2.1
1	A	242	ARG	2.1
1	A	575	HIS	2.1
1	A	467	ARG	2.0

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](#)

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	PEG	A	601	7/7	0.85	0.19	25,46,59,71	0

6.5 Other polymers [i](#)

There are no such residues in this entry.