



Full wwPDB X-ray Structure Validation Report ⓘ

May 25, 2020 – 08:41 am BST

PDB ID : 3QSK
Title : 5 Histidine Variant of the anti-RNase A VHH in Complex with RNase A
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Deposited on : 2011-02-21
Resolution : 1.75 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 following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
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

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 2026 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 Ribonuclease pancreatic.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	119	921	558	165	186	12	0	1	0

- Molecule 2 is a protein called Engineered 5 Histidine anti-RNase A Camelid VHH Antibody Domain Variant.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	B	123	923	571	167	180	5	0	1	0

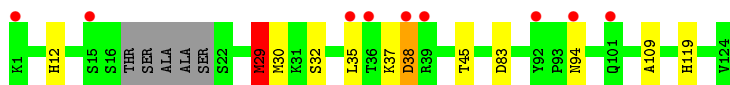
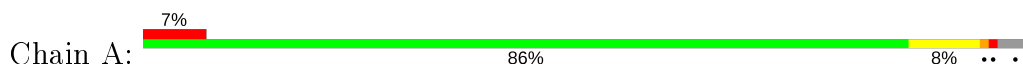
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	72	Total 72	O 72	0	0
3	B	110	Total 110	O 110	0	0

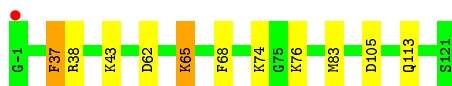
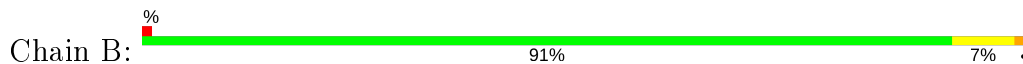
3 Residue-property plots [i](#)

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.

- Molecule 1: Ribonuclease pancreatic



- Molecule 2: Engineered 5 Histidine anti-RNase A Camelid VHH Antibody Domain Variant



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	40.79Å 54.68Å 48.79Å 90.00° 109.24° 90.00°	Depositor
Resolution (Å)	20.00 – 1.75 19.25 – 1.75	Depositor EDS
% Data completeness (in resolution range)	99.8 (20.00-1.75) 99.8 (19.25-1.75)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	10.42 (at 1.75Å)	Xtrriage
Refinement program	REFMAC 5.5.0102	Depositor
R, R_{free}	0.184 , 0.246 0.188 , 0.247	Depositor DCC
R_{free} test set	1023 reflections (4.99%)	wwPDB-VP
Wilson B-factor (Å ²)	16.5	Xtrriage
Anisotropy	0.144	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 43.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	2026	wwPDB-VP
Average B, all atoms (Å ²)	21.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 10.09% 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](#)

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	0.96	0/939	0.89	1/1265 (0.1%)
2	B	1.20	1/947 (0.1%)	1.03	3/1277 (0.2%)
All	All	1.09	1/1886 (0.1%)	0.96	4/2542 (0.2%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	37	PHE	CD1-CE1	5.03	1.49	1.39

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	68	PHE	CB-CG-CD2	-5.83	116.72	120.80
2	B	83	MET	CG-SD-CE	-5.58	91.28	100.20
1	A	29	MET	CG-SD-CE	5.17	108.47	100.20
2	B	38	ARG	NE-CZ-NH2	-5.02	117.79	120.30

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 clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	921	0	871	10	0
2	B	923	0	875	4	0
3	A	72	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	110	0	0	2	0
All	All	2026	0	1746	14	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 4.

All (14) 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:30:MET:HE3	1:A:35:LEU:HB2	1.51	0.92
1:A:30:MET:CE	1:A:35:LEU:HB2	2.16	0.76
1:A:30:MET:HE3	1:A:35:LEU:CB	2.28	0.62
1:A:30:MET:CE	1:A:35:LEU:CB	2.78	0.60
1:A:12:HIS:HD2	1:A:45:THR:O	1.85	0.60
1:A:109:ALA:HB3	1:A:119:HIS:HB3	1.90	0.53
2:B:113:GLN:HG2	3:B:1061:HOH:O	2.07	0.53
1:A:37:LYS:O	1:A:38:ASP:HB3	2.09	0.52
1:A:37:LYS:O	1:A:38:ASP:CB	2.58	0.51
2:B:37:PHE:CE2	2:B:105:ASP:HB3	2.46	0.51
2:B:43:LYS:HD3	3:B:1087:HOH:O	2.12	0.50
1:A:29:MET:HE3	1:A:32:SER:HB2	1.95	0.48
1:A:29:MET:HA	1:A:29:MET:CE	2.44	0.47
2:B:62:ASP:O	2:B:65:LYS:HG2	2.14	0.47

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	116/124 (94%)	111 (96%)	4 (3%)	1 (1%)	17 5
2	B	122/123 (99%)	121 (99%)	1 (1%)	0	100 100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
All	All	238/247 (96%)	232 (98%)	5 (2%)	1 (0%)	34 17

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	38	ASP

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	106/109 (97%)	103 (97%)	3 (3%)	43 20
2	B	93/92 (101%)	90 (97%)	3 (3%)	39 16
All	All	199/201 (99%)	193 (97%)	6 (3%)	41 18

All (6) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	29	MET
1	A	83	ASP
1	A	94	ASN
2	B	65	LYS
2	B	74	LYS
2	B	76	LYS

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

Mol	Chain	Res	Type
1	A	12	HIS

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 carbohydrates 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](#)

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	119/124 (95%)	0.48	9 (7%) 13 18	10, 25, 43, 50	2 (1%)
2	B	123/123 (100%)	-0.17	1 (0%) 86 90	8, 15, 26, 37	1 (0%)
All	All	242/247 (97%)	0.15	10 (4%) 37 44	8, 18, 38, 50	3 (1%)

All (10) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	92	TYR	4.5
1	A	38	ASP	3.3
1	A	39	ARG	3.2
1	A	101	GLN	3.0
2	B	-1	GLY	2.8
1	A	15	SER	2.7
1	A	35	LEU	2.4
1	A	94	ASN	2.4
1	A	36	THR	2.3
1	A	1	LYS	2.1

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](#)

There are no ligands in this entry.

6.5 Other polymers

There are no such residues in this entry.