



Full wwPDB X-ray Structure Validation Report ⓘ

Apr 16, 2025 – 12:07 PM JST

PDB ID : 9KEP / pdb_00009kep
Title : RABV-G-PHD-FD/NM57-scFv
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Deposited on : 2024-11-05
Resolution : 2.89 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
Xtriage (Phenix)	:	2.0rc1
EDS	:	3.0
Percentile statistics	:	20231227.v01 (using entries in the PDB archive December 27th 2023)
CCP4	:	9.0.006 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.42

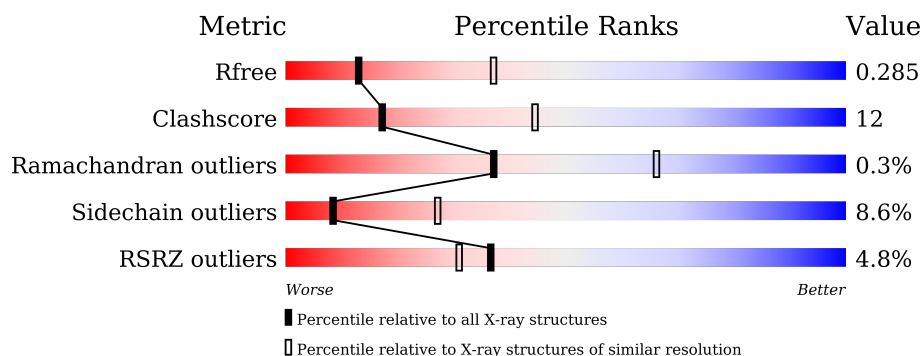
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.89 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}	164625	2335 (2.90-2.90)
Clashscore	180529	2564 (2.90-2.90)
Ramachandran outliers	177936	2514 (2.90-2.90)
Sidechain outliers	177891	2516 (2.90-2.90)
RSRZ outliers	164620	2337 (2.90-2.90)

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	116	<div> <div>2%</div> <div>76%</div> <div>21%</div> <div>• •</div> </div>
2	B	243	<div> <div>7%</div> <div>39%</div> <div>21%</div> <div>•</div> <div>37%</div> </div>
3	D	127	<div> <div>%</div> <div>68%</div> <div>27%</div> <div>• •</div> </div>

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 2984 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 NM57-scFv light chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	113	Total	C	N	O	S	0	0	0
			824	512	138	170	4			

- Molecule 2 is a protein called RABV-G-PHD-FD.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	154	Total	C	N	O	S	0	0	0
			1194	747	209	224	14			

- Molecule 3 is a protein called NM57-scFv heavy chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	D	125	Total	C	N	O	S	0	0	0
			966	611	163	188	4			

4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	114.33Å 49.48Å 100.86Å 90.00° 117.62° 90.00°	Depositor
Resolution (Å)	29.79 – 2.89 29.79 – 2.89	Depositor EDS
% Data completeness (in resolution range)	99.9 (29.79-2.89) 99.8 (29.79-2.89)	Depositor EDS
R_{merge}	0.14	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.91 (at 2.90Å)	Xtriage
Refinement program	PHENIX (1.20.1_4487: ???)	Depositor
R, R_{free}	0.233 , 0.281 0.234 , 0.285	Depositor DCC
R_{free} test set	525 reflections (4.53%)	wwPDB-VP
Wilson B-factor (Å ²)	69.9	Xtriage
Anisotropy	0.480	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 63.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	2984	wwPDB-VP
Average B, all atoms (Å ²)	87.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.12% 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

5.1 Standard geometry

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	0/843	0.55	0/1145
2	B	0.47	1/1217 (0.1%)	0.91	7/1637 (0.4%)
3	D	0.28	0/990	0.57	1/1345 (0.1%)
All	All	0.36	1/3050 (0.0%)	0.72	8/4127 (0.2%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
2	B	0	1
3	D	0	1
All	All	0	2

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	89	PRO	N-CD	12.48	1.65	1.47

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	200	ALA	CB-CA-C	-15.08	87.47	110.10
2	B	201	SER	N-CA-CB	-10.09	95.36	110.50
2	B	89	PRO	CA-N-CD	-8.62	99.43	111.50
2	B	89	PRO	N-CA-CB	6.55	111.16	103.30
2	B	201	SER	CB-CA-C	6.48	122.42	110.10
2	B	88	ARG	N-CA-C	6.15	127.61	111.00
3	D	86	LEU	CB-CG-CD1	-5.84	101.07	111.00
2	B	215	LEU	N-CA-CB	-5.30	99.81	110.40

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	B	198	LYS	Peptide
3	D	12	LYS	Peptide

5.2 Too-close contacts

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	824	0	786	14	0
2	B	1194	0	1139	41	0
3	D	966	0	916	20	0
All	All	2984	0	2841	72	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 12.

All (72) 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)
2:B:88:ARG:HB3	2:B:89:PRO:HD3	1.57	0.84
2:B:37:ASN:CB	2:B:200:ALA:O	2.36	0.72
2:B:37:ASN:HB3	2:B:200:ALA:O	1.89	0.72
3:D:91:THR:HG23	3:D:124:THR:HA	1.71	0.72
2:B:208:GLY:HA2	2:B:219:LEU:HD13	1.74	0.68
2:B:171:THR:OG1	2:B:176:THR:OG1	2.13	0.66
1:A:48:LEU:HD21	1:A:51:TYR:HB3	1.76	0.66
3:D:12:LYS:HG3	3:D:18:VAL:HG22	1.78	0.65
1:A:52:ASP:OD1	1:A:53:ALA:N	2.32	0.63
1:A:1:GLN:NE2	3:D:46:GLU:OE2	2.32	0.61
2:B:30:VAL:HG23	2:B:31:GLU:H	1.67	0.59
2:B:41:PHE:HZ	2:B:243:MET:SD	2.28	0.57
2:B:199:ARG:HB2	2:B:209:PHE:HA	1.86	0.57
2:B:201:SER:HB3	2:B:205:LYS:HE3	1.87	0.56
2:B:37:ASN:HB2	2:B:200:ALA:O	2.05	0.55
2:B:43:TYR:CZ	2:B:195:SER:HB2	2.42	0.54
2:B:59:PHE:CE1	2:B:179:MET:HB2	2.43	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:62:ASP:OD1	1:A:62:ASP:N	2.36	0.54
2:B:41:PHE:HB3	2:B:196:ARG:O	2.08	0.54
3:D:115:ASP:HB3	3:D:116:PRO:HD3	1.90	0.53
2:B:233:LEU:HD11	2:B:245:THR:HG22	1.91	0.53
3:D:73:ASP:OD2	3:D:76:THR:OG1	2.25	0.52
2:B:197:GLY:HA3	2:B:210:VAL:O	2.11	0.51
1:A:26:SER:HB2	1:A:93:TYR:O	2.10	0.51
2:B:159:CYS:SG	2:B:160:SER:N	2.84	0.51
3:D:13:LYS:HB2	3:D:16:SER:OG	2.11	0.50
3:D:60:TYR:CE2	3:D:70:ILE:HD12	2.45	0.50
3:D:87:ARG:HB2	3:D:89:ASP:OD1	2.12	0.50
2:B:233:LEU:N	2:B:241:VAL:O	2.45	0.50
1:A:18:VAL:HG21	1:A:109:LEU:HD13	1.94	0.49
3:D:29:PHE:CD2	3:D:100:ASN:HB3	2.47	0.49
3:D:85:SER:O	3:D:85:SER:OG	2.28	0.49
2:B:95:ARG:NH2	2:B:180:PRO:O	2.45	0.49
2:B:90:THR:O	2:B:94:CYS:HB2	2.13	0.48
2:B:223:CYS:SG	2:B:234:ARG:NE	2.78	0.48
1:A:38:TYR:OH	3:D:112:GLY:O	2.32	0.48
2:B:61:CYS:O	2:B:176:THR:HA	2.13	0.48
2:B:240:TRP:CH2	2:B:242:ALA:HB2	2.49	0.47
1:A:83:GLU:OE2	1:A:83:GLU:N	2.42	0.47
2:B:88:ARG:HA	2:B:175:TYR:HA	1.96	0.47
2:B:44:MET:HG2	2:B:240:TRP:HZ3	1.80	0.47
2:B:59:PHE:HE1	2:B:94:CYS:SG	2.39	0.46
1:A:52:ASP:HB3	1:A:55:LYS:HE2	1.97	0.46
2:B:43:TYR:CE1	2:B:195:SER:HB2	2.51	0.46
3:D:62:GLN:O	3:D:65:GLN:HB3	2.15	0.46
2:B:44:MET:HB2	2:B:44:MET:HE2	1.81	0.46
2:B:233:LEU:CD1	2:B:243:MET:HB2	2.46	0.45
2:B:192:PHE:HZ	2:B:227:LEU:HD12	1.80	0.45
2:B:199:ARG:HB2	2:B:208:GLY:O	2.16	0.45
2:B:88:ARG:CZ	2:B:89:PRO:HG3	2.47	0.45
1:A:7:PRO:O	1:A:107:THR:OG1	2.27	0.44
2:B:141:ASP:O	2:B:150:HIS:N	2.48	0.44
3:D:2:VAL:HG22	3:D:26:GLY:O	2.17	0.44
1:A:34:PHE:CZ	3:D:113:TRP:HB2	2.52	0.43
2:B:224:ARG:N	2:B:250:LYS:O	2.40	0.43
3:D:18:VAL:CG2	3:D:86:LEU:HD11	2.47	0.43
2:B:46:LEU:HD12	2:B:46:LEU:HA	1.80	0.43
2:B:202:ASN:HB3	2:B:205:LYS:HE2	2.01	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:D:65:GLN:O	3:D:67:ARG:HG3	2.19	0.43
3:D:46:GLU:OE2	3:D:63:ARG:HD3	2.19	0.42
2:B:88:ARG:HB3	2:B:89:PRO:CD	2.39	0.42
2:B:235:LEU:HB2	2:B:237:ASP:OD1	2.19	0.42
3:D:67:ARG:NH2	3:D:90:ASP:OD2	2.51	0.42
1:A:63:ARG:HB3	1:A:78:SER:O	2.20	0.42
2:B:43:TYR:CE2	2:B:45:GLU:HB2	2.56	0.41
1:A:51:TYR:CE2	1:A:55:LYS:HD2	2.55	0.41
2:B:223:CYS:SG	2:B:234:ARG:NH2	2.91	0.41
3:D:81:MET:O	3:D:81:MET:HG3	2.20	0.41
3:D:18:VAL:HG23	3:D:86:LEU:HD21	2.03	0.41
2:B:237:ASP:OD1	2:B:238:GLY:N	2.54	0.40
1:A:96:ASP:HB2	1:A:97:TYR:CD1	2.57	0.40
2:B:192:PHE:CZ	2:B:227:LEU:HD12	2.55	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	111/116 (96%)	108 (97%)	3 (3%)	0	100	100
2	B	138/243 (57%)	128 (93%)	9 (6%)	1 (1%)	19	49
3	D	123/127 (97%)	113 (92%)	10 (8%)	0	100	100
All	All	372/486 (76%)	349 (94%)	22 (6%)	1 (0%)	37	66

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	89	PRO

5.3.2 Protein sidechains ⓘ

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	90/93 (97%)	84 (93%)	6 (7%)	13	39
2	B	133/210 (63%)	124 (93%)	9 (7%)	13	38
3	D	102/104 (98%)	89 (87%)	13 (13%)	3	11
All	All	325/407 (80%)	297 (91%)	28 (9%)	8	27

All (28) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	28	ASP
1	A	65	SER
1	A	67	SER
1	A	96	ASP
1	A	108	LYS
1	A	110	THR
2	B	32	ASP
2	B	61	CYS
2	B	86	HIS
2	B	136	SER
2	B	158	LYS
2	B	175	TYR
2	B	199	ARG
2	B	228	CYS
2	B	244	GLN
3	D	16	SER
3	D	21	SER
3	D	22	CYS
3	D	33	THR
3	D	35	ASN
3	D	59	ASN
3	D	63	ARG
3	D	75	SER
3	D	77	SER
3	D	87	ARG
3	D	96	CYS

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Mol	Chain	Res	Type
3	D	103	ASN
3	D	109	TYR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
2	B	172	ASN
2	B	202	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 oligosaccharides 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

6.1 Protein, DNA and RNA chains

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	113/116 (97%)	0.02	2 (1%) 67 61	39, 56, 80, 109	0
2	B	154/243 (63%)	1.01	16 (10%) 13 12	61, 119, 150, 170	0
3	D	125/127 (98%)	0.12	1 (0%) 82 78	48, 73, 106, 122	0
All	All	392/486 (80%)	0.44	19 (4%) 36 31	39, 80, 143, 170	0

All (19) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	90	THR	3.7
2	B	63	GLY	3.5
2	B	89	PRO	3.5
2	B	133	ILE	3.4
2	B	200	ALA	3.3
2	B	35	CYS	3.1
2	B	211	ASP	2.9
2	B	209	PHE	2.9
3	D	115	ASP	2.8
2	B	149	LEU	2.6
2	B	101	LEU	2.6
2	B	86	HIS	2.5
2	B	37	ASN	2.4
2	B	167	THR	2.3
1	A	54	THR	2.2
2	B	46	LEU	2.2
1	A	51	TYR	2.1
2	B	245	THR	2.0
2	B	186	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](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.