



Full wwPDB X-ray Structure Validation Report ⓘ

May 25, 2020 – 10:31 pm BST

PDB ID : 6ICC
Title : The NZ-1 Fab complexed with the PDZ tandem fragment of A. aeolicus S2P homolog with the PA12 tag inserted between the residues 181 and 186
Authors : Tamura, R.; Oi, R.; Kaneko, M.K.; Kato, Y.; Nogi, T.
Deposited on : 2018-09-05
Resolution : 2.00 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
Mogul : 1.8.5 (274361), CSD as541be (2020)
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

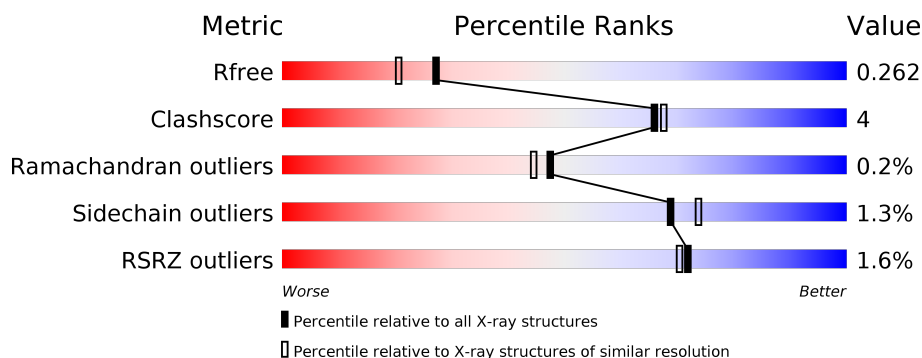
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.00 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	8085 (2.00-2.00)
Clashscore	141614	9178 (2.00-2.00)
Ramachandran outliers	138981	9054 (2.00-2.00)
Sidechain outliers	138945	9053 (2.00-2.00)
RSRZ outliers	127900	7900 (2.00-2.00)

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 on 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	188	<div> <div>2%</div> <div> <div></div> <div>90%</div> <div>8%</div> <div></div> </div> </div>
2	H	219	<div> <div>3%</div> <div> <div></div> <div>91%</div> <div>9%</div> </div> </div>
3	L	214	<div> <div></div> <div> <div></div> <div>93%</div> <div>7%</div> </div> </div>

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 5046 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 PDZ tandem fragment with PA tag.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	185	Total	C	N	O	S	0	5	0
			1470	955	246	267	2			

There are 14 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	113	GLY	-	expression tag	UNP O67776
A	114	SER	-	expression tag	UNP O67776
A	181A	GLY	-	see sequence details	UNP O67776
A	181B	VAL	-	see sequence details	UNP O67776
A	181C	ALA	-	see sequence details	UNP O67776
A	181D	MET	-	see sequence details	UNP O67776
A	181E	PRO	-	see sequence details	UNP O67776
A	181F	GLY	-	see sequence details	UNP O67776
A	181G	ALA	-	see sequence details	UNP O67776
A	181H	GLU	-	see sequence details	UNP O67776
A	181I	ASP	-	see sequence details	UNP O67776
A	181J	ASP	-	see sequence details	UNP O67776
A	181K	VAL	-	see sequence details	UNP O67776
A	181L	VAL	-	see sequence details	UNP O67776

- Molecule 2 is a protein called Heavy chain of antigen binding fragment, Fab of NZ-1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	H	219	Total	C	N	O	S	0	1	0
			1650	1040	275	326	9			

- Molecule 3 is a protein called Light chain of antigen binding fragment, Fab of NZ-1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	L	214	Total	C	N	O	S	0	4	0
			1675	1036	293	340	6			

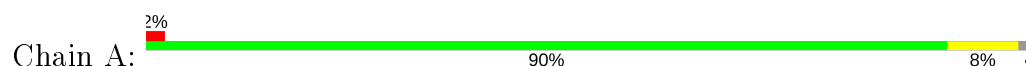
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	62	Total 62	O 62	0	0
4	H	87	Total 87	O 87	0	0
4	L	102	Total 102	O 102	0	0

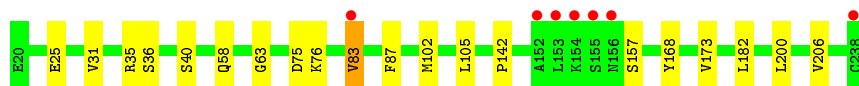
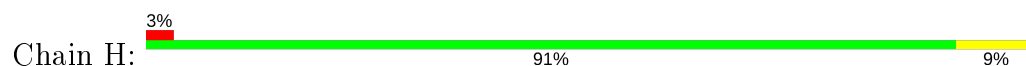
3 Residue-property plots

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.

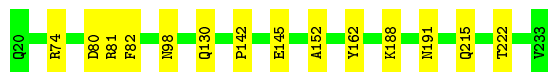
- Molecule 1: PDZ tandem fragment with PA tag



- Molecule 2: Heavy chain of antigen binding fragment, Fab of NZ-1



- Molecule 3: Light chain of antigen binding fragment, Fab of NZ-1



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	49.94Å 84.29Å 185.91Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	48.23 – 2.00 48.23 – 2.00	Depositor EDS
% Data completeness (in resolution range)	100.0 (48.23-2.00) 100.0 (48.23-2.00)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.42 (at 2.00Å)	Xtriage
Refinement program	REFMAC 5.8.0232	Depositor
R, R_{free}	0.213 , 0.260 0.214 , 0.262	Depositor DCC
R_{free} test set	2701 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	41.5	Xtriage
Anisotropy	0.073	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 41.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	5046	wwPDB-VP
Average B, all atoms (Å ²)	47.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.22% 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: SNN, PCA

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	0.29	0/1493	0.60	0/2011
2	H	0.31	0/1690	0.64	0/2303
3	L	0.30	0/1703	0.62	0/2319
All	All	0.30	0/4886	0.62	0/6633

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	1470	0	1580	14	0
2	H	1650	0	1610	17	0
3	L	1675	0	1607	7	0
4	A	62	0	0	3	0
4	H	87	0	0	1	0
4	L	102	0	0	1	0
All	All	5046	0	4797	38	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 (38) 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:176[A]:THR:HG23	1:A:187:HIS:HB2	1.31	1.10
1:A:176[A]:THR:HG23	1:A:187:HIS:CB	2.04	0.87
1:A:176[A]:THR:CG2	1:A:187:HIS:HB2	2.14	0.73
2:H:182:LEU:HD21	2:H:206:VAL:HG11	1.73	0.70
2:H:102:MET:HE2	2:H:105:LEU:HD21	1.74	0.68
1:A:149[B]:ILE:HD11	1:A:177:LEU:HD13	1.75	0.67
3:L:188:LYS:HE2	3:L:191:ASN:HA	1.78	0.66
3:L:130:GLN:HB2	3:L:162:TYR:CE1	2.32	0.64
1:A:149[B]:ILE:CD1	1:A:177:LEU:HD13	2.27	0.63
2:H:76:LYS:HA	4:H:326:HOH:O	1.99	0.62
1:A:176[A]:THR:CG2	1:A:187:HIS:CB	2.77	0.61
2:H:102:MET:CE	2:H:105:LEU:HD21	2.33	0.57
2:H:182:LEU:HD21	2:H:206:VAL:CG1	2.36	0.56
1:A:187:HIS:HD2	4:A:345:HOH:O	1.90	0.55
3:L:215:GLN:HG3	3:L:222:THR:CG2	2.37	0.55
2:H:200:LEU:C	2:H:200:LEU:HD12	2.28	0.54
1:A:223[A]:VAL:HG12	1:A:270:LYS:HD2	1.89	0.54
2:H:31:VAL:HG22	2:H:35:ARG:HB2	1.90	0.53
1:A:187:HIS:CD2	4:A:345:HOH:O	2.63	0.50
2:H:36:SER:HA	2:H:102:MET:O	2.12	0.49
3:L:81:ARG:HB2	3:L:98[B]:ASN:O	2.12	0.48
2:H:25:GLU:HA	2:H:40[A]:SER:O	2.16	0.46
2:H:83:VAL:HG21	2:H:87:PHE:HB2	1.98	0.46
3:L:74[A]:ARG:HD3	3:L:82:PHE:O	2.15	0.46
2:H:58:GLN:HA	2:H:63:GLY:O	2.17	0.45
1:A:196:GLN:HG2	4:A:337:HOH:O	2.17	0.45
2:H:83:VAL:O	2:H:83:VAL:HG22	2.17	0.45
2:H:31:VAL:HG11	2:H:105:LEU:HD13	2.00	0.44
1:A:176[A]:THR:HG22	1:A:178:PHE:CE1	2.53	0.43
2:H:25:GLU:HA	2:H:40[B]:SER:O	2.18	0.43
3:L:145:GLU:HB2	4:L:382:HOH:O	2.18	0.43
2:H:173:VAL:CG2	2:H:200:LEU:HD21	2.48	0.43
2:H:142:PRO:HB3	2:H:168:TYR:HB3	2.00	0.43
1:A:140:ILE:HD11	1:A:177:LEU:CD2	2.50	0.41
1:A:160:LEU:HD21	1:A:203:ILE:HD11	2.02	0.41
2:H:31:VAL:CG2	2:H:35:ARG:HB2	2.50	0.41
3:L:142:PRO:HG3	3:L:152:ALA:HB1	2.02	0.41

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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	184/188 (98%)	182 (99%)	1 (0%)	1 (0%)	29	23
2	H	218/219 (100%)	214 (98%)	4 (2%)	0	100	100
3	L	216/214 (101%)	210 (97%)	6 (3%)	0	100	100
All	All	618/621 (100%)	606 (98%)	11 (2%)	1 (0%)	47	44

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	206	LEU

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	162/159 (102%)	159 (98%)	3 (2%)	57	61
2	H	184/183 (100%)	181 (98%)	3 (2%)	62	67
3	L	191/187 (102%)	190 (100%)	1 (0%)	88	92
All	All	537/529 (102%)	530 (99%)	7 (1%)	69	74

All (7) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	170	ASP
1	A	238	LYS

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Mol	Chain	Res	Type
1	A	258	LYS
2	H	75	ASP
2	H	83	VAL
2	H	157	SER
3	L	80	ASP

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

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

2 non-standard protein/DNA/RNA residues 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	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	PCA	L	20	3	7,8,9	0.51	0	9,10,12	0.94	0
1	SNN	A	150	1	7,8,8	1.85	2 (28%)	7,11,11	2.86	3 (42%)

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
3	PCA	L	20	3	-	0/0/11/13	0/1/1/1
1	SNN	A	150	1	-	-	0/1/1/1

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	150	SNN	C2-N1	-3.73	1.32	1.37
1	A	150	SNN	C5-N1	-3.02	1.32	1.37

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	150	SNN	C3-C2-N1	5.33	111.42	107.30
1	A	150	SNN	O2-C2-C3	-4.34	123.05	126.18
1	A	150	SNN	O5-C5-C4	-2.60	122.96	126.39

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

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	184/188 (97%)	0.24	3 (1%) 72 70	30, 48, 69, 85	0
2	H	219/219 (100%)	0.28	7 (3%) 47 46	31, 46, 71, 95	0
3	L	213/214 (99%)	0.04	0 100 100	27, 43, 60, 77	0
All	All	616/621 (99%)	0.19	10 (1%) 72 70	27, 45, 68, 95	0

All (10) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	H	154	LYS	6.1
2	H	155	SER	4.6
2	H	238	CYS	4.5
2	H	153	LEU	3.6
2	H	83	VAL	3.5
2	H	156	ASN	3.3
1	A	266	LYS	3.2
1	A	187	HIS	3.0
2	H	152	ALA	3.0
1	A	262	LEU	2.2

6.2 Non-standard residues in protein, DNA, RNA chains [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
3	PCA	L	20	8/9	0.72	0.21	77,84,86,88	0
1	SNN	A	150	8/8	0.95	0.12	35,38,41,43	0

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 [i](#)

There are no such residues in this entry.