



wwPDB X-ray Structure Validation Summary Report ⓘ

Feb 1, 2016 – 01:40 AM GMT

PDB ID : 2DW2
Title : Crystal structure of VAP2 from Crotalus atrox venom (Form 2-5 crystal)
Authors : Takeda, S.; Igarashi, T.; Araki, S.
Deposited on : 2006-08-02
Resolution : 2.70 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
<http://wwpdb.org/validation/2016/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 : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

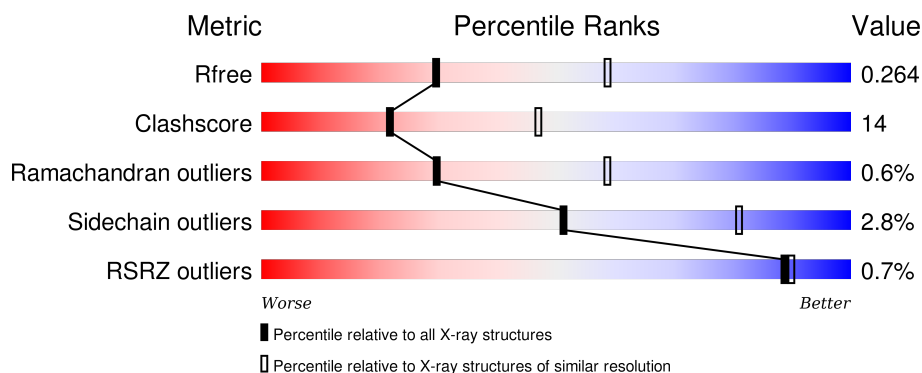
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.70 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}	91344	2103 (2.70-2.70)
Clashscore	102246	2422 (2.70-2.70)
Ramachandran outliers	100387	2382 (2.70-2.70)
Sidechain outliers	100360	2382 (2.70-2.70)
RSRZ outliers	91569	2107 (2.70-2.70)

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. 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	419	<div> <div></div> <div>74%24%..</div> </div>
1	B	419	<div> <div></div> <div>70%27%..</div> </div>

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 6823 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 Catrocollastatin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	415	Total	C	N	O	S	0	0	0
			3219	1990	545	637	47			
1	B	415	Total	C	N	O	S	0	0	0
			3219	1990	545	637	47			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	203	VAL	PHE	SEE REMARK 999	UNP Q90282
B	203	VAL	PHE	SEE REMARK 999	UNP Q90282

- Molecule 2 is a polymer of unknown type called SUGAR (9-MER).

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	9	Total	C	N	O	0	0
			113	64	5	44		
2	B	9	Total	C	N	O	0	0
			113	64	5	44		

- Molecule 3 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	1	Total	Zn	0	0
			1	1		
3	A	1	Total	Zn	0	0
			1	1		

- Molecule 4 is CALCIUM ION (three-letter code: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	B	3	Total	Ca	0	0
			3	3		

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	3	Total	Ca	0	0
			3	3		

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	85	Total	O	0	0
			85	85		
5	B	66	Total	O	0	0
			66	66		

4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	220.74Å 79.47Å 58.69Å 90.00° 91.73° 90.00°	Depositor
Resolution (Å)	50.00 – 2.70 46.38 – 2.70	Depositor EDS
% Data completeness (in resolution range)	95.8 (50.00-2.70) 95.6 (46.38-2.70)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	5.31 (at 2.69Å)	Xtriage
Refinement program	CNS 1.0	Depositor
R, R_{free}	0.199 , 0.260 0.214 , 0.264	Depositor DCC
R_{free} test set	1334 reflections (4.96%)	DCC
Wilson B-factor (Å ²)	26.0	Xtriage
Anisotropy	0.373	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 46.1	EDS
Estimated twinning fraction	0.094 for -h,-k,l	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Outliers	0 of 26911 reflections	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	6823	wwPDB-VP
Average B, all atoms (Å ²)	25.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.49% 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, 0.375 respectively for untwinned datasets, and 0.333, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: ZN, BMA, NAG, CA, FUC, MAN

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	0/3291	0.55	0/4453
1	B	0.28	0/3291	0.55	0/4453
All	All	0.28	0/6582	0.55	0/8906

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

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	3219	0	2985	82	0
1	B	3219	0	2985	98	0
2	A	113	0	97	5	0
2	B	113	0	97	1	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
4	A	3	0	0	0	0
4	B	3	0	0	0	0
5	A	85	0	0	4	0
5	B	66	0	0	5	0
All	All	6823	0	6164	180	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 14.

The worst 5 of 180 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:529:ASN:HD21	1:A:557:ARG:HB3	1.36	0.91
1:B:326:VAL:HG13	1:B:361:ILE:HD11	1.63	0.81
1:B:529:ASN:HD21	1:B:557:ARG:HB3	1.47	0.80
1:B:454:LYS:HE3	1:B:457:THR:HG21	1.69	0.72
1:B:454:LYS:CD	1:B:454:LYS:H	2.02	0.72

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	413/419 (99%)	380 (92%)	31 (8%)	2 (0%)	34	63
1	B	413/419 (99%)	381 (92%)	29 (7%)	3 (1%)	26	55
All	All	826/838 (99%)	761 (92%)	60 (7%)	5 (1%)	30	59

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	430	ALA
1	A	568	ASN
1	B	277	ASP
1	B	196	PRO
1	A	566	GLY

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	361/365 (99%)	355 (98%)	6 (2%)	68	90
1	B	361/365 (99%)	347 (96%)	14 (4%)	39	70
All	All	722/730 (99%)	702 (97%)	20 (3%)	51	81

5 of 20 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	388	CYS
1	B	391	ASN
1	B	523	ASP
1	B	360	GLU
1	B	385	ASN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 27 such sidechains are listed below:

Mol	Chain	Res	Type
1	B	268	ASN
1	B	385	ASN
1	B	563	ASN
1	B	288	GLN
1	A	485	HIS

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates

18 carbohydrates 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$
2	NAG	A	801	1,2	14,14,15	0.55	0	15,19,21	0.83	1 (6%)
2	NAG	A	802	2	14,14,15	0.51	0	15,19,21	0.87	1 (6%)
2	BMA	A	803	2	11,11,12	0.52	0	14,15,17	0.58	0
2	NAG	A	804	2	14,14,15	0.46	0	15,19,21	0.77	1 (6%)
2	MAN	A	805	2	11,11,12	0.92	1 (9%)	14,15,17	1.70	2 (14%)
2	NAG	A	806	2	14,14,15	0.64	0	15,19,21	0.75	1 (6%)
2	MAN	A	807	2	11,11,12	0.64	0	14,15,17	0.63	0
2	NAG	A	808	2	14,14,15	0.56	0	15,19,21	0.70	1 (6%)
2	FUC	A	809	2	10,10,11	0.52	0	14,14,16	0.59	0
2	NAG	B	801	1,2	14,14,15	0.53	0	15,19,21	0.86	1 (6%)
2	NAG	B	802	2	14,14,15	0.54	0	15,19,21	0.76	0
2	BMA	B	803	2	11,11,12	0.52	0	14,15,17	0.63	0
2	NAG	B	804	2	14,14,15	0.44	0	15,19,21	0.82	1 (6%)
2	MAN	B	805	2	11,11,12	0.96	1 (9%)	14,15,17	1.53	1 (7%)
2	NAG	B	806	2	14,14,15	0.67	0	15,19,21	0.78	1 (6%)
2	MAN	B	807	2	11,11,12	0.38	0	14,15,17	0.75	1 (7%)
2	NAG	B	808	2	14,14,15	0.48	0	15,19,21	0.65	0
2	FUC	B	809	2	10,10,11	0.64	0	14,14,16	0.54	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	NAG	A	801	1,2	-	0/6/23/26	0/1/1/1
2	NAG	A	802	2	-	0/6/23/26	0/1/1/1
2	BMA	A	803	2	-	0/2/19/22	0/1/1/1
2	NAG	A	804	2	-	0/6/23/26	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	MAN	A	805	2	-	0/2/19/22	0/1/1/1
2	NAG	A	806	2	-	0/6/23/26	0/1/1/1
2	MAN	A	807	2	-	0/2/19/22	0/1/1/1
2	NAG	A	808	2	-	0/6/23/26	0/1/1/1
2	FUC	A	809	2	-	0/0/17/20	0/1/1/1
2	NAG	B	801	1,2	-	0/6/23/26	0/1/1/1
2	NAG	B	802	2	-	0/6/23/26	0/1/1/1
2	BMA	B	803	2	-	0/2/19/22	0/1/1/1
2	NAG	B	804	2	-	0/6/23/26	0/1/1/1
2	MAN	B	805	2	-	0/2/19/22	0/1/1/1
2	NAG	B	806	2	-	0/6/23/26	0/1/1/1
2	MAN	B	807	2	-	0/2/19/22	0/1/1/1
2	NAG	B	808	2	-	0/6/23/26	0/1/1/1
2	FUC	B	809	2	-	0/0/17/20	0/1/1/1

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	805	MAN	C2-C3	2.45	1.55	1.52
2	A	805	MAN	C2-C3	2.51	1.56	1.52

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

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	804	NAG	C2-N2-C7	-2.57	119.73	123.04
2	A	801	NAG	C2-N2-C7	-2.53	119.79	123.04
2	A	804	NAG	C2-N2-C7	-2.45	119.89	123.04
2	A	806	NAG	C2-N2-C7	-2.20	120.21	123.04
2	B	806	NAG	C2-N2-C7	-2.14	120.30	123.04

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

6 monomers are involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	804	NAG	1	0
2	A	805	MAN	2	0
2	A	806	NAG	3	0
2	A	807	MAN	1	0

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Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	808	NAG	1	0
2	B	806	NAG	1	0

5.6 Ligand geometry [i](#)

Of 8 ligands modelled in this entry, 8 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

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	415/419 (99%)	-0.17	3 (0%) 89 90	6, 22, 43, 66	0
1	B	415/419 (99%)	-0.12	3 (0%) 89 90	8, 24, 46, 61	0
All	All	830/838 (99%)	-0.14	6 (0%) 89 90	6, 23, 44, 66	0

The worst 5 of 6 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	568	ASN	3.8
1	A	567	GLN	3.4
1	B	542	ASN	3.0
1	B	424	GLN	2.8
1	A	352	ASP	2.6

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](#)

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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors(Å ²)	Q<0.9
2	NAG	A	806	14/15	0.91	0.19	1.29	14,31,33,34	0
2	NAG	B	804	14/15	0.93	0.16	0.49	24,31,34,36	0
2	NAG	B	801	14/15	0.92	0.17	0.21	29,34,41,46	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
2	NAG	B	806	14/15	0.89	0.17	0.07	28,35,38,39	0
2	NAG	A	804	14/15	0.94	0.16	0.06	30,35,40,40	0
2	NAG	A	801	14/15	0.94	0.15	-0.01	26,31,35,37	0
2	FUC	A	809	10/11	0.93	0.23	-	23,30,34,39	0
2	NAG	A	802	14/15	0.93	0.14	-	24,32,35,37	0
2	FUC	B	809	10/11	0.92	0.21	-	26,30,34,36	0
2	BMA	A	803	11/12	0.94	0.15	-	29,35,38,39	0
2	MAN	A	807	11/12	0.87	0.17	-	32,38,45,46	0
2	NAG	B	808	14/15	0.86	0.26	-	42,47,48,51	0
2	NAG	B	802	14/15	0.93	0.18	-	24,39,45,46	0
2	BMA	B	803	11/12	0.96	0.20	-	34,40,42,45	0
2	MAN	A	805	11/12	0.90	0.24	-	34,37,39,48	0
2	MAN	B	805	11/12	0.89	0.24	-	32,40,43,45	0
2	MAN	B	807	11/12	0.93	0.18	-	33,36,39,40	0
2	NAG	A	808	14/15	0.91	0.30	-	42,45,48,49	0

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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors(\AA^2)	Q<0.9
4	CA	B	711	1/1	0.96	0.12	-0.50	23,23,23,23	0
4	CA	A	703	1/1	0.98	0.13	-1.64	23,23,23,23	0
4	CA	B	713	1/1	0.98	0.12	-1.70	20,20,20,20	0
4	CA	A	702	1/1	0.97	0.06	-2.77	20,20,20,20	0
4	CA	A	701	1/1	0.99	0.06	-2.84	21,21,21,21	0
4	CA	B	712	1/1	0.99	0.06	-3.14	22,22,22,22	0
3	ZN	A	700	1/1	1.00	0.10	-3.56	18,18,18,18	0
3	ZN	B	700	1/1	1.00	0.12	-	17,17,17,17	0

6.5 Other polymers [i](#)

There are no such residues in this entry.