



Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 10:48 PM GMT

PDB ID : 1VAP
Title : THE MONOMERIC ASP49 SECRETORY PHOSPHOLIPASE A2 FROM
THE VENOM OF AGKISTRIDON PISCIVORUS PISCIVORUS
Authors : Scott, D.L.
Deposited on : 1996-11-15
Resolution : 1.60 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
<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) : **NOT EXECUTED**
EDS : **NOT EXECUTED**
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

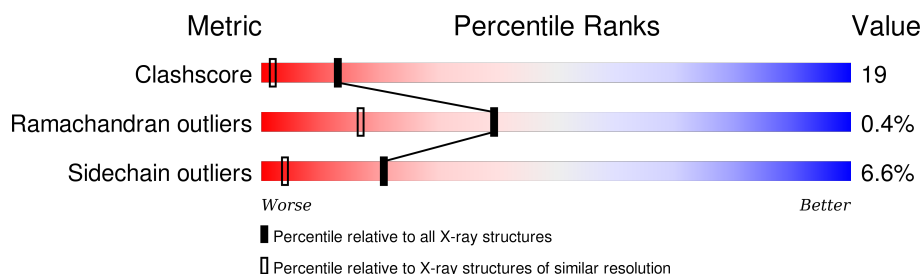
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.60 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))
Clashscore	102246	2732 (1.60-1.60)
Ramachandran outliers	100387	2654 (1.60-1.60)
Sidechain outliers	100360	2653 (1.60-1.60)

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.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	123	 66% 30% ..
1	B	123	 66% 31% ..

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 2081 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 PHOSPHOLIPASE A2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	123	Total	C	N	O	S	0	0	0
			971	608	165	181	17			
1	B	123	Total	C	N	O	S	0	0	0
			971	608	165	181	17			

- Molecule 2 is water.

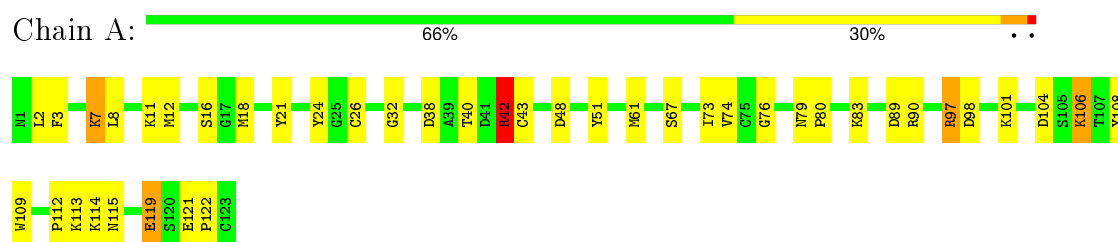
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	81	Total	O	0	0
			81	81		
2	B	58	Total	O	0	0
			58	58		

3 Residue-property plots

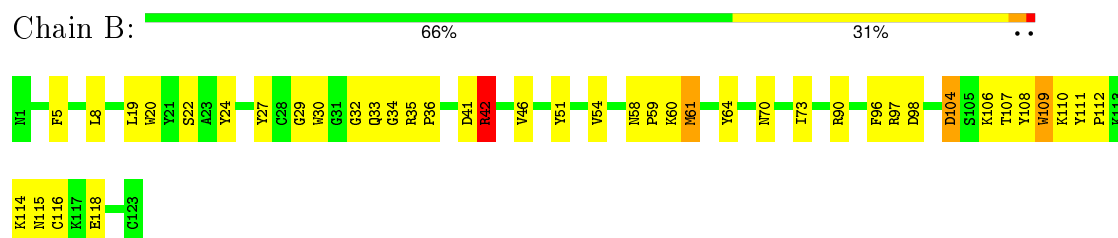
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according 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.

Note EDS was not executed.

• Molecule 1: PHOSPHOLIPASE A2



• Molecule 1: PHOSPHOLIPASE A2



4 Data and refinement statistics

Xtriage (Phenix) and EDS were not executed - this section will therefore be incomplete.

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	34.71Å 39.89Å 42.85Å 94.17° 79.10° 106.27°	Depositor
Resolution (Å)	10.00 – 1.60	Depositor
% Data completeness (in resolution range)	(Not available) (10.00-1.60)	Depositor
R_{merge}	0.04	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	PROFFT, X-PLOR	Depositor
R, R_{free}	0.197 , (Not available)	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	2081	wwPDB-VP
Average B, all atoms (Å ²)	22.0	wwPDB-VP

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	0.82	0/995	1.47	13/1335 (1.0%)
1	B	0.82	0/995	1.44	10/1335 (0.7%)
All	All	0.82	0/1990	1.45	23/2670 (0.9%)

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
1	A	0	2
1	B	0	1
All	All	0	3

There are no bond length outliers.

All (23) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	90	ARG	NE-CZ-NH2	-13.30	113.65	120.30
1	A	89	ASP	CB-CG-OD2	-9.84	109.45	118.30
1	B	64	TYR	CB-CG-CD2	-8.47	115.92	121.00
1	A	108	TYR	CB-CG-CD2	-8.24	116.06	121.00
1	B	98	ASP	CB-CG-OD2	-8.13	110.98	118.30
1	A	24	TYR	CB-CG-CD1	7.60	125.56	121.00
1	A	24	TYR	CB-CG-CD2	-7.45	116.53	121.00
1	B	90	ARG	NH1-CZ-NH2	7.39	127.53	119.40
1	A	48	ASP	CB-CG-OD1	6.99	124.59	118.30
1	A	98	ASP	CB-CG-OD2	-6.83	112.15	118.30
1	A	90	ARG	NE-CZ-NH2	-6.67	116.97	120.30
1	B	5	PHE	CB-CG-CD2	-6.52	116.24	120.80
1	B	97	ARG	NE-CZ-NH1	6.26	123.43	120.30
1	B	96	PHE	CB-CG-CD2	-6.05	116.56	120.80

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	7	LYS	N-CA-CB	5.62	120.71	110.60
1	B	104	ASP	CB-CG-OD2	-5.60	113.26	118.30
1	A	3	PHE	N-CA-CB	5.53	120.56	110.60
1	A	21	TYR	CB-CG-CD2	5.53	124.31	121.00
1	A	108	TYR	CB-CG-CD1	5.52	124.31	121.00
1	A	38	ASP	O-C-N	5.48	131.47	122.70
1	A	51	TYR	CB-CG-CD2	-5.36	117.78	121.00
1	B	42	ARG	O-C-N	5.27	131.13	122.70
1	B	96	PHE	CB-CG-CD1	5.08	124.36	120.80

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	42	ARG	Sidechain
1	A	97	ARG	Sidechain
1	B	46	VAL	Mainchain

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	971	0	919	40	0
1	B	971	0	919	32	0
2	A	81	0	0	7	1
2	B	58	0	0	1	1
All	All	2081	0	1838	72	1

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 19.

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)
1:A:113:LYS:HE2	2:A:285:HOH:O	1.49	1.10

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:106:LYS:H	1:A:106:LYS:HD3	0.90	1.02
1:A:106:LYS:CD	1:A:106:LYS:H	1.73	1.02
1:A:106:LYS:N	1:A:106:LYS:HD3	1.74	0.97
1:B:70:ASN:HB2	2:B:263:HOH:O	1.65	0.95
1:A:112:PRO:HG2	1:A:115:ASN:OD1	1.70	0.91
1:A:104:ASP:OD1	1:A:106:LYS:HE2	1.71	0.90
1:A:104:ASP:OD1	1:A:106:LYS:CE	2.19	0.90
1:B:112:PRO:HB2	1:B:114:LYS:HG2	1.54	0.89
1:A:42:ARG:HH11	1:A:42:ARG:HG3	1.38	0.87
1:A:113:LYS:CE	2:A:285:HOH:O	2.11	0.84
1:B:54:VAL:HG21	1:B:59:PRO:HG3	1.61	0.81
1:B:51:TYR:O	1:B:54:VAL:HG22	1.81	0.80
1:B:54:VAL:CG2	1:B:59:PRO:HG3	2.11	0.80
1:A:42:ARG:NH1	1:A:42:ARG:HG3	1.96	0.79
1:A:114:LYS:CG	1:A:115:ASN:ND2	2.47	0.77
1:B:22:SER:O	1:B:29:GLY:HA3	1.85	0.76
1:A:114:LYS:HG3	1:A:115:ASN:ND2	2.02	0.75
1:A:97:ARG:HD3	2:A:306:HOH:O	1.86	0.73
1:A:112:PRO:CG	1:A:115:ASN:OD1	2.36	0.72
1:A:104:ASP:OD1	1:A:106:LYS:HE3	1.90	0.71
1:B:111:TYR:CD1	1:B:112:PRO:HD2	2.27	0.68
1:B:112:PRO:HG2	1:B:115:ASN:OD1	1.94	0.67
1:B:54:VAL:HG21	1:B:59:PRO:CG	2.24	0.66
1:A:114:LYS:HG2	1:A:115:ASN:ND2	2.10	0.66
1:B:58:ASN:ND2	1:B:61:MET:HB2	2.11	0.66
1:B:111:TYR:CG	1:B:112:PRO:HD2	2.34	0.62
1:A:114:LYS:HG3	1:A:115:ASN:HD22	1.64	0.61
1:B:58:ASN:HD21	1:B:61:MET:HE3	1.65	0.61
1:B:35:ARG:HG3	1:B:118:GLU:HG2	1.82	0.61
1:B:70:ASN:O	1:B:70:ASN:CG	2.37	0.61
1:A:12:MET:HE3	1:A:97:ARG:HD2	1.84	0.59
1:A:26:CYS:O	1:A:32:GLY:HA2	2.03	0.59
1:A:97:ARG:CD	2:A:239:HOH:O	2.52	0.57
1:A:97:ARG:HD3	2:A:239:HOH:O	2.05	0.56
1:B:58:ASN:HD21	1:B:61:MET:CE	2.17	0.56
1:A:115:ASN:ND2	1:A:115:ASN:N	2.52	0.56
1:B:24:TYR:HD2	1:B:109:TRP:HD1	1.52	0.56
1:B:19:LEU:O	1:B:109:TRP:CZ2	2.58	0.56
1:A:114:LYS:CG	1:A:115:ASN:HD21	2.19	0.53
1:B:33:GLN:HG3	1:B:116:CYS:O	2.09	0.53
1:A:12:MET:CE	1:A:97:ARG:HD2	2.39	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:109:TRP:CD1	2:A:322:HOH:O	2.54	0.52
1:B:22:SER:O	1:B:29:GLY:CA	2.57	0.50
1:A:112:PRO:HG2	1:A:115:ASN:CG	2.30	0.50
1:A:40:THR:O	1:A:43:CYS:HB2	2.12	0.50
1:A:115:ASN:N	1:A:115:ASN:HD22	2.09	0.49
1:A:42:ARG:NH1	1:A:42:ARG:CG	2.71	0.49
1:A:112:PRO:C	1:A:114:LYS:H	2.15	0.49
1:A:7:LYS:HE2	2:A:281:HOH:O	2.12	0.48
1:A:8:LEU:HD22	1:A:73:ILE:HD13	1.94	0.48
1:B:36:PRO:HA	1:B:41:ASP:OD2	2.15	0.47
1:B:34:GLY:HA2	1:B:118:GLU:O	2.13	0.47
1:B:42:ARG:HB2	1:B:42:ARG:CZ	2.45	0.46
1:B:35:ARG:HG3	1:B:118:GLU:CG	2.45	0.46
1:A:11:LYS:HB2	1:A:11:LYS:HE2	1.33	0.45
1:A:67:SER:OG	1:A:74:VAL:CG1	2.64	0.45
1:B:35:ARG:HG2	1:B:35:ARG:HH11	1.81	0.45
1:B:104:ASP:OD2	1:B:107:THR:HG23	2.17	0.44
1:B:58:ASN:ND2	1:B:61:MET:HE3	2.32	0.44
1:B:20:TRP:O	1:B:109:TRP:NE1	2.45	0.44
1:A:18:MET:CA	1:A:18:MET:HE2	2.48	0.43
1:B:107:THR:OG1	1:B:108:TYR:CD2	2.70	0.43
1:A:67:SER:O	1:A:74:VAL:HG12	2.18	0.43
1:B:104:ASP:OD2	1:B:107:THR:CG2	2.67	0.43
1:A:112:PRO:C	1:A:114:LYS:N	2.72	0.42
1:B:8:LEU:HG	1:B:73:ILE:HD13	2.01	0.41
1:A:121:GLU:HG3	1:A:122:PRO:HD2	2.03	0.41
1:B:112:PRO:C	1:B:114:LYS:N	2.73	0.41
1:A:76:GLY:O	1:A:83:LYS:HG2	2.21	0.40
1:A:79:ASN:HA	1:A:80:PRO:HD3	1.79	0.40
1:B:27:TYR:CD1	1:B:32:GLY:HA3	2.57	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:227:HOH:O	2:B:323:HOH:O[1_554]	2.04	0.16

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	121/123 (98%)	114 (94%)	6 (5%)	1 (1%)	24	6
1	B	121/123 (98%)	117 (97%)	4 (3%)	0	100	100
All	All	242/246 (98%)	231 (96%)	10 (4%)	1 (0%)	39	17

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	119	GLU

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	106/106 (100%)	99 (93%)	7 (7%)	21	4
1	B	106/106 (100%)	99 (93%)	7 (7%)	21	4
All	All	212/212 (100%)	198 (93%)	14 (7%)	21	4

All (14) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	2	LEU
1	A	16	SER
1	A	42	ARG
1	A	61	MET
1	A	101	LYS

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Mol	Chain	Res	Type
1	A	106	LYS
1	A	119	GLU
1	B	30	TRP
1	B	42	ARG
1	B	60	LYS
1	B	61	MET
1	B	106	LYS
1	B	109	TRP
1	B	110	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	33	GLN
1	A	70	ASN
1	A	84	GLN
1	A	115	ASN
1	B	58	ASN

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 ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry ⓘ

There are no ligands in this entry.

5.7 Other polymers ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues ⓘ

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

EDS was not executed - this section will therefore be empty.

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

EDS was not executed - this section will therefore be empty.

6.3 Carbohydrates ⓘ

EDS was not executed - this section will therefore be empty.

6.4 Ligands ⓘ

EDS was not executed - this section will therefore be empty.

6.5 Other polymers ⓘ

EDS was not executed - this section will therefore be empty.