



# Full wwPDB X-ray Structure Validation Report

Feb 27, 2014 – 07:57 PM GMT

PDB ID : 1VWB  
Title : STREPTAVIDIN-CYCLO-AC-[CHPQFC]-NH<sub>2</sub>,PH 11.8  
Authors : Katz, B.A.; Cass, R.T.  
Deposited on : 1997-03-03  
Resolution : 1.82 Å(reported)

This is a full wwPDB validation report for a publicly released PDB entry.  
We welcome your comments at [validation@mail.wwpdb.org](mailto:validation@mail.wwpdb.org)  
A user guide is available at <http://wwpdb.org/ValidationPDFNotes.html>

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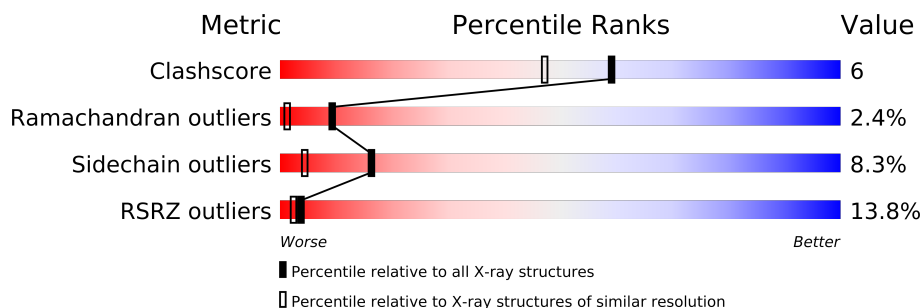
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467  
Mogul : 1.15 2013  
Xtriage (Phenix) : dev-1323  
EDS : stable22639  
Percentile statistics : 21963  
Refmac : 5.8.0049  
CCP4 : 6.3.0 (Settle)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)  
Validation Pipeline (wwPDB-VP) : stable22683

# 1 Overall quality at a glance

The reported resolution of this entry is 1.82 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	79885	5140 (1.84-1.80)
Ramachandran outliers	78287	5077 (1.84-1.80)
Sidechain outliers	78261	5077 (1.84-1.80)
RSRZ outliers	66119	4103 (1.84-1.80)

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  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	B	123	
2	P	8	

## 2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 1426 atoms, of which 352 are hydrogens and 0 are deuterium.

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 STREPTAVIDIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	B	123	Total	C	H	N	O	12	8	0
			1204	612	236	167	189			

- Molecule 2 is a protein called PEPTIDE LIGAND CONTAINING HPQ.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	P	8	Total	C	H	N	O	0	0	1
			63	33	10	10	8			

- Molecule 3 is water.

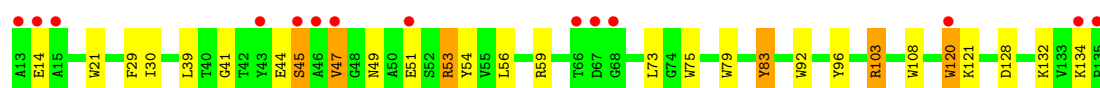
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	B	53	Total	H	O	0	0
			159	106	53		

### 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.

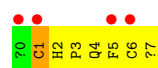
- Molecule 1: STREPTAVIDIN

Chain B:



- Molecule 2: PEPTIDE LIGAND CONTAINING HPQ

Chain P:



## 4 Data and refinement statistics

Property	Value	Source
Space group	I 41 2 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	59.25Å 59.25Å 178.35Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	7.50 – 1.82 28.11 – 1.73	Depositor EDS
% Data completeness (in resolution range)	75.9 (7.50-1.82) 67.7 (28.11-1.73)	Depositor EDS
$R_{merge}$	0.07	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.36 (at 1.73Å)	Xtriage
Refinement program	X-PLOR	Depositor
R, $R_{free}$	0.183 , 0.239 0.234 , (Not available)	Depositor DCC
$R_{free}$ test set	No test flags present.	DCC
Wilson B-factor (Å <sup>2</sup> )	17.2	Xtriage
Anisotropy	0.176	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.37 , 90.4	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 11592 reflections	Xtriage
$F_o, F_c$ correlation	0.90	EDS
Total number of atoms	1426	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	19.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.79% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

## 5 Model quality

### 5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: ACE, NH2

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	B	0.91	0/1029	1.62	27/1404 (1.9%)
2	P	0.98	0/52	1.53	1/70 (1.4%)
All	All	0.91	0/1081	1.61	28/1474 (1.9%)

There are no bond length outliers.

All (28) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	108	TRP	CD1-CG-CD2	9.90	114.22	106.30
1	B	120	TRP	CD1-CG-CD2	8.90	113.42	106.30
1	B	108	TRP	CE2-CD2-CG	-8.67	100.36	107.30
1	B	92	TRP	CD1-CG-CD2	7.95	112.66	106.30
1	B	120	TRP	CE2-CD2-CG	-7.78	101.08	107.30
1	B	75	TRP	CD1-CG-CD2	7.77	112.52	106.30
1	B	75	TRP	CE2-CD2-CG	-7.60	101.22	107.30
1	B	21	TRP	CD1-CG-CD2	7.32	112.16	106.30
1	B	92	TRP	CE2-CD2-CG	-7.20	101.54	107.30
1	B	120	TRP	CG-CD2-CE3	7.18	140.36	133.90
1	B	120	TRP	CB-CG-CD1	-7.14	117.72	127.00
1	B	79	TRP	CD1-CG-CD2	7.01	111.91	106.30
1	B	79	TRP	CE2-CD2-CG	-6.54	102.07	107.30
1	B	83	TYR	CB-CG-CD2	-6.38	117.17	121.00
1	B	21	TRP	CE2-CD2-CG	-6.38	102.20	107.30
1	B	92	TRP	CG-CD2-CE3	6.14	139.43	133.90
2	P	6	CYS	CA-CB-SG	-6.13	102.97	114.00
1	B	92	TRP	CB-CG-CD1	-5.86	119.38	127.00
1	B	120	TRP	CG-CD1-NE1	-5.83	104.27	110.10
1	B	59	ARG	NE-CZ-NH1	5.79	123.20	120.30
1	B	108	TRP	CB-CG-CD1	-5.59	119.73	127.00
1	B	108	TRP	CG-CD2-CE3	5.55	138.90	133.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	54	TYR	CB-CG-CD2	-5.37	117.78	121.00
1	B	108	TRP	CG-CD1-NE1	-5.32	104.78	110.10
1	B	96	TYR	CB-CG-CD2	-5.28	117.83	121.00
1	B	92	TRP	CG-CD1-NE1	-5.21	104.89	110.10
1	B	103	ARG	CA-CB-CG	5.16	124.75	113.40
1	B	53	ARG	NE-CZ-NH1	5.14	122.87	120.30

There are no chirality outliers.

There are no planarity outliers.

## 5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	B	968	236	694	10	0
2	P	53	10	34	4	0
3	B	53	106	0	1	0
All	All	1074	352	728	12	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 6.

All (12) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:120:TRP:CZ3	1:B:121:LYS:HB3	2.40	0.57
1:B:120:TRP:CE3	1:B:121:LYS:HB3	2.44	0.53
1:B:45:SER:HB2	2:P:7:NH2:N	2.24	0.53
1:B:44:GLU:HG3	1:B:53:ARG:HG2	1.91	0.53
1:B:39:LEU:HD11	1:B:73[B]:LEU:HD12	1.91	0.51
1:B:51:GLU:HA	1:B:83:TYR:CD2	2.49	0.47
1:B:45:SER:HB3	2:P:3:PRO:O	2.15	0.46
1:B:44:GLU:HB2	3:B:613:HOH:H2	1.81	0.46
1:B:29[B]:PHE:HD2	1:B:56:LEU:HD21	1.82	0.45
2:P:4:GLN:HB3	2:P:5:PHE:CE2	2.55	0.42
2:P:2:HIS:HB3	2:P:5:PHE:HD2	1.86	0.41

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:30:ILE:O	1:B:41:GLY:HA3	2.21	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	B	130/123 (106%)	124 (95%)	4 (3%)	2 (2%)	15	3
2	P	6/8 (75%)	5 (83%)	0	1 (17%)	0	0
All	All	136/131 (104%)	129 (95%)	4 (3%)	3 (2%)	9	1

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	14	GLU
2	P	1	CYS
1	B	47	VAL

### 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	B	99/90 (110%)	92 (93%)	7 (7%)	21	6
2	P	6/6 (100%)	5 (83%)	1 (17%)	3	0
All	All	105/96 (109%)	97 (92%)	8 (8%)	16	5

All (8) residues with a non-rotameric sidechain are listed below:



Mol	Chain	Res	Type
1	B	45	SER
1	B	47	VAL
1	B	49	ASN
1	B	103	ARG
1	B	128	ASP
1	B	132	LYS
1	B	134	LYS
2	P	1	CYS

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 chains 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 ⓘ

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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	B	122/123 (99%)	0.75	14 (11%) 5 4	10, 17, 50, 78	22 (18%)
2	P	8/8 (100%)	2.41	4 (50%) 0 0	27, 39, 48, 50	4 (50%)
All	All	130/131 (99%)	0.85	18 (13%) 4 2	10, 19, 50, 78	26 (20%)

All (18) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	47	VAL	16.2
1	B	13	ALA	13.0
2	P	1	CYS	9.4
1	B	135	PRO	7.2
1	B	14	GLU	4.9
1	B	15	ALA	4.9
2	P	5	PHE	4.0
2	P	6	CYS	3.6
1	B	66	THR	3.1
1	B	68	GLY	3.0
1	B	45	SER	3.0
1	B	67	ASP	2.8
1	B	120	TRP	2.8
1	B	46	ALA	2.6
1	B	43[A]	TYR	2.6
2	P	0	ACE	2.4
1	B	134	LYS	2.4
1	B	51	GLU	2.1

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

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

### 6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

### 6.4 Ligands ⓘ

There are no ligands in this entry.

### 6.5 Other polymers ⓘ

There are no such residues in this entry.