



Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 10:12 PM GMT

PDB ID : 1SLE
Title : STREPTAVIDIN, PH 5.0, BOUND TO CYCLIC PEPTIDE AC-CHPQGPPC-NH2
Authors : Katz, B.A.
Deposited on : 1995-03-10
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
<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

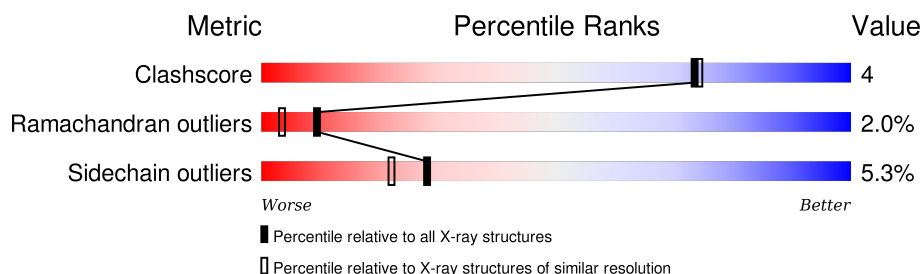
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 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))
Clashscore	102246	7340 (2.00-2.00)
Ramachandran outliers	100387	7248 (2.00-2.00)
Sidechain outliers	100360	7247 (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. 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	B	135	
1	D	135	
2	M	10	
2	P	10	

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 2756 atoms, of which 705 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 STREPTAVIDIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	B	121	Total	C	H	N	O	3	0	0
			1113	562	212	156	183			
1	D	121	Total	C	H	N	O	0	0	0
			1113	562	212	156	183			

- Molecule 2 is a protein called AC-CHPQGPPC-NH2.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
2	M	10	Total	C	H	N	O	S	0	0	2
			70	36	11	12	9	2			
2	P	10	Total	C	H	N	O	S	0	0	1
			70	36	10	12	10	2			



- Molecule 3 is water.



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	B	61	Total	H	O	0	0
			183	122	61		
3	D	64	Total	H	O	0	0
			192	128	64		
3	P	5	Total	H	O	0	0
			15	10	5		

Note EDS was not executed.

- Chain B:
-
- 70% 16% 10%
- | Amino Acid | Category |
|------------|-----------|
| ASP | Basic |
| PRO | Non-polar |
| SER | Polar |
| LYS | Basic |
| ASP | Basic |
| SER | Polar |
| LYS | Basic |
| ALA | Non-polar |
| GLN | Polar |
| VAL | Non-polar |
| SER | Polar |
| ALA | Non-polar |
| A13 | Non-polar |
| E14 | Acidic |
| W21 | Aromatic |
| Y22 | Aromatic |
| T28 | Non-polar |
| A35 | Non-polar |
| L39 | Non-polar |
| E44 | Acidic |
| S45 | Polar |
| A46 | Non-polar |
| V47 | Non-polar |
| G48 | Polar |
| M49 | Non-polar |
| A50 | Non-polar |
| E51 | Acidic |
| S52 | Polar |
| Y60 | Aromatic |
| T66 | Non-polar |
| D67 | Acidic |
| G70 | Polar |
| L73 | Non-polar |
| G74 | Polar |
| W75 | Aromatic |
| W79 | Aromatic |
| R84 | Basic |
| W92 | Aromatic |
| Q95 | Polar |
| Q107 | Polar |
| W108 | Aromatic |
| T114 | Non-polar |
| W120 | Aromatic |

- Chain D:

- Chain M:  80% 20%
- 
- 70% C1 H2 P3 79%

- Chain P: 
- 

4 Data and refinement statistics

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

Property	Value	Source
Space group	I 2 2 2	Depositor
Cell constants a, b, c, α , β , γ	97.65Å 107.61Å 49.46Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	7.50 – 2.00	Depositor
% Data completeness (in resolution range)	(Not available) (7.50-2.00)	Depositor
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	X-PLOR	Depositor
R, R_{free}	0.191 , 0.233	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	2756	wwPDB-VP
Average B, all atoms (Å ²)	16.0	wwPDB-VP

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	0.96	0/923	1.65	24/1264 (1.9%)
1	D	0.95	0/923	1.69	25/1264 (2.0%)
2	M	0.83	0/60	1.28	0/82
2	P	0.72	0/60	1.40	0/83
All	All	0.94	0/1966	1.65	49/2693 (1.8%)

There are no bond length outliers.

All (49) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	120	TRP	CD1-CG-CD2	8.27	112.91	106.30
1	D	92	TRP	CD1-CG-CD2	8.04	112.73	106.30
1	B	75	TRP	CE2-CD2-CG	-7.86	101.01	107.30
1	D	21	TRP	CD1-CG-CD2	7.79	112.53	106.30
1	D	51	GLU	CA-C-N	-7.73	100.20	117.20
1	B	79	TRP	CD1-CG-CD2	7.59	112.37	106.30
1	B	75	TRP	CD1-CG-CD2	7.53	112.32	106.30
1	B	108	TRP	CE2-CD2-CG	-7.47	101.33	107.30
1	B	75	TRP	CG-CD2-CE3	7.33	140.49	133.90
1	B	108	TRP	CD1-CG-CD2	7.28	112.12	106.30
1	D	21	TRP	CE2-CD2-CG	-7.25	101.50	107.30
1	D	92	TRP	CE2-CD2-CG	-7.13	101.59	107.30
1	B	92	TRP	CD1-CG-CD2	6.87	111.79	106.30
1	D	75	TRP	CD1-CG-CD2	6.80	111.74	106.30
1	B	79	TRP	CE2-CD2-CG	-6.80	101.86	107.30
1	D	79	TRP	CE2-CD2-CG	-6.79	101.87	107.30
1	D	120	TRP	CE2-CD2-CG	-6.70	101.94	107.30
1	D	108	TRP	CE2-CD2-CG	-6.65	101.98	107.30
1	D	79	TRP	CD1-CG-CD2	6.64	111.61	106.30
1	B	120	TRP	CE2-CD2-CG	-6.62	102.01	107.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	75	TRP	CE2-CD2-CG	-6.61	102.01	107.30
1	D	43	TYR	CB-CG-CD2	-6.50	117.10	121.00
1	B	75	TRP	CB-CG-CD1	-6.47	118.58	127.00
1	D	96	TYR	CB-CG-CD2	-6.46	117.13	121.00
1	D	120	TRP	CD1-CG-CD2	6.43	111.45	106.30
1	D	92	TRP	CB-CG-CD1	-6.24	118.89	127.00
1	D	121	LYS	CA-CB-CG	6.14	126.91	113.40
1	B	92	TRP	CE2-CD2-CG	-6.07	102.44	107.30
1	D	108	TRP	CD1-CG-CD2	6.00	111.10	106.30
1	B	21	TRP	CE2-CD2-CG	-5.95	102.54	107.30
1	B	67	ASP	N-CA-C	-5.80	95.34	111.00
1	D	54	TYR	CB-CG-CD2	-5.80	117.52	121.00
1	D	92	TRP	CG-CD2-CE3	5.67	139.00	133.90
1	B	21	TRP	CD1-CG-CD2	5.66	110.83	106.30
1	B	120	TRP	CG-CD1-NE1	-5.60	104.50	110.10
1	D	31	VAL	CB-CA-C	-5.55	100.85	111.40
1	B	73	LEU	CA-CB-CG	5.42	127.77	115.30
1	B	39	LEU	CA-CB-CG	5.39	127.71	115.30
1	B	45	SER	N-CA-CB	-5.39	102.42	110.50
1	B	66	THR	CA-C-N	-5.21	105.74	117.20
1	D	103	ARG	CB-CA-C	-5.19	100.02	110.40
1	B	79	TRP	CG-CD1-NE1	-5.17	104.93	110.10
1	D	128	ASP	CB-CG-OD1	5.12	122.91	118.30
1	B	84	ARG	NE-CZ-NH1	5.12	122.86	120.30
1	D	77	VAL	CG1-CB-CG2	-5.11	102.72	110.90
1	D	75	TRP	CG-CD2-CE3	5.08	138.48	133.90
1	B	108	TRP	CB-CG-CD1	-5.04	120.45	127.00
1	B	107	GLN	CA-CB-CG	-5.02	102.36	113.40
1	D	120	TRP	CG-CD2-CE3	5.00	138.40	133.90

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	B	901	212	839	8	1
1	D	901	212	839	8	0
2	M	59	11	47	1	0
2	P	60	10	50	1	0
3	B	61	122	0	0	0
3	D	64	128	0	1	1
3	P	5	10	0	0	0
All	All	2051	705	1775	15	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 4.

All (15) 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:B:95:GLN:HE22	1:D:114:THR:H	1.31	0.79
1:B:114:THR:H	1:D:95:GLN:HE22	1.36	0.73
1:D:44:GLU:HB3	1:D:53:ARG:HG2	1.83	0.60
2:P:0:ACE:H1	2:P:1:CYS:SG	2.52	0.49
1:D:67:ASP:HB3	3:D:709:HOH:O	2.13	0.48
1:B:70:GLY:HA3	1:B:95:GLN:HE21	1.81	0.45
1:B:14:GLU:HB3	1:B:60:TYR:OH	2.16	0.45
1:D:14:GLU:HB3	1:D:60:TYR:OH	2.15	0.45
1:D:70:GLY:HA3	1:D:95:GLN:HE21	1.80	0.45
1:B:22:TYR:HE1	1:B:28:THR:HG1	1.65	0.44
1:B:28:THR:HB	1:B:44:GLU:HB2	1.99	0.43
1:B:45:SER:HB2	2:M:3:PRO:O	2.20	0.42
1:D:14:GLU:HG3	1:D:15:ALA:H	1.85	0.41
1:D:49:ASN:OD1	1:D:50:ALA:N	2.54	0.41
1:B:49:ASN:HB3	1:B:52:SER:OG	2.20	0.41

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)
1:B:35:ALA:O	3:D:921:HOH:H1[6_555]	1.60	0.00

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	119/135 (88%)	111 (93%)	5 (4%)	3 (2%)	7	2
1	D	119/135 (88%)	111 (93%)	7 (6%)	1 (1%)	24	15
2	M	8/10 (80%)	5 (62%)	2 (25%)	1 (12%)	0	0
2	P	8/10 (80%)	8 (100%)	0	0	100	100
All	All	254/290 (88%)	235 (92%)	14 (6%)	5 (2%)	9	3

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	46	ALA
1	B	47	VAL
1	B	67	ASP
2	M	1	CYS
1	D	50	ALA

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	88/100 (88%)	84 (96%)	4 (4%)	34	29
1	D	88/100 (88%)	82 (93%)	6 (7%)	20	13
2	M	7/7 (100%)	7 (100%)	0	100	100
2	P	7/7 (100%)	7 (100%)	0	100	100
All	All	190/214 (89%)	180 (95%)	10 (5%)	28	22

All (10) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	39	LEU
1	B	51	GLU
1	B	52	SER
1	B	73	LEU
1	D	14	GLU
1	D	31	VAL
1	D	36	ASP
1	D	66	THR
1	D	103	ARG
1	D	121	LYS

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

Mol	Chain	Res	Type
1	B	95	GLN
1	D	95	GLN

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 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 ⓘ

There are no chain breaks in this entry.

6 Fit of model and data [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

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

6.3 Carbohydrates [i](#)

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

6.4 Ligands [i](#)

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

6.5 Other polymers [i](#)

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