



Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 10:13 PM GMT

PDB ID : 1SLG
Title : STREPTAVIDIN, PH 5.6, BOUND TO PEPTIDE FCHPQNT
Authors : Katz, B.A.
Deposited on : 1995-03-10
Resolution : 1.76 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

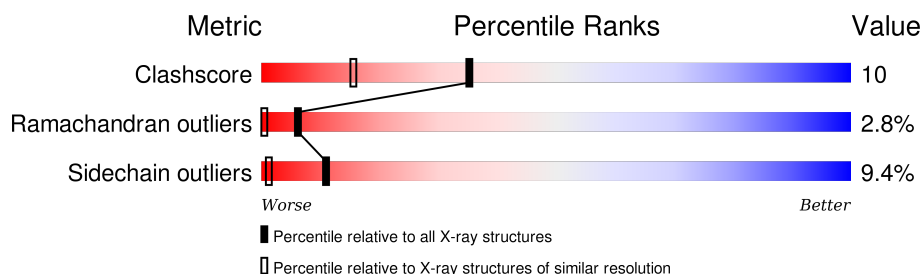
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.76 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	1730 (1.76-1.76)
Ramachandran outliers	100387	1711 (1.76-1.76)
Sidechain outliers	100360	1711 (1.76-1.76)

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	 62% 26% 9%
1	D	135	 64% 19% 5% 10%
2	M	7	 43% 43% 14%
2	P	7	 29% 43% 29%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 2913 atoms, of which 784 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	123	Total	C	H	N	O	0	3	0
			1152	588	218	159	187			
1	D	121	Total	C	H	N	O	11	4	0
			1137	575	219	158	185			

- Molecule 2 is a protein called FCHPQNT.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	M	6	Total	C	H	N	O	0	0	0
			59	27	11	10	11			
2	P	7	Total	C	H	N	O	0	1	0
			88	45	18	12	13			

- Molecule 3 is water.

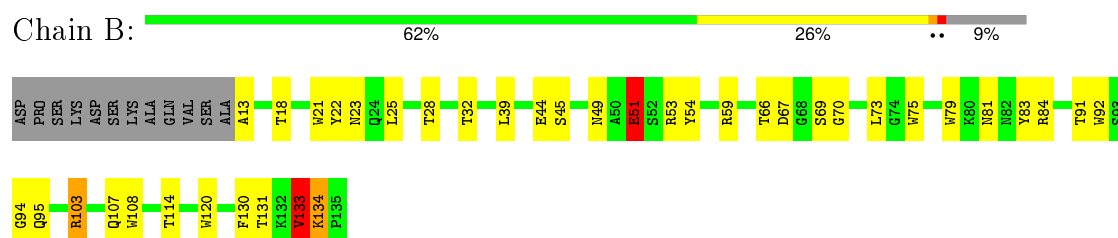
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	B	75	Total	H	O	0	0
			225	150	75		
3	D	80	Total	H	O	0	0
			240	160	80		
3	M	3	Total	H	O	0	0
			9	6	3		
3	P	1	Total	H	O	0	0
			3	2	1		

3 Residue-property plots [i](#)

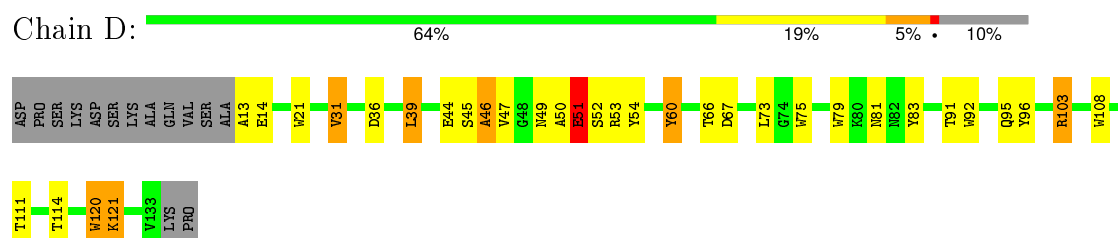
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according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

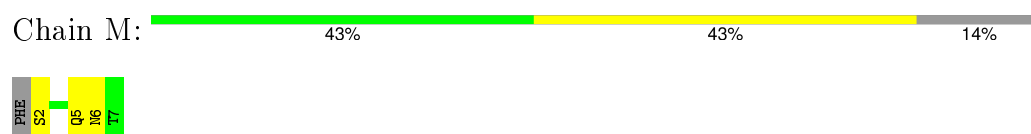
• Molecule 1: STREPTAVIDIN



• Molecule 1: STREPTAVIDIN



• Molecule 2: FCHPQNT



• Molecule 2: FCHPQNT



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	96.24 Å 105.89 Å 48.02 Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	7.50 – 1.76	Depositor
% Data completeness (in resolution range)	(Not available) (7.50-1.76)	Depositor
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	X-PLOR	Depositor
R, R_{free}	0.187 , 0.222	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	2913	wwPDB-VP
Average B, all atoms (Å ²)	17.0	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	B	0.98	0/970	1.83	25/1327 (1.9%)
1	D	0.95	0/956	1.78	28/1309 (2.1%)
2	M	0.91	0/49	1.47	0/65
2	P	1.41	2/73 (2.7%)	2.08	3/97 (3.1%)
All	All	0.99	2/2048 (0.1%)	1.81	56/2798 (2.0%)

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	B	0	2
1	D	0	1
All	All	0	3

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	P	2	SER	CA-CB	6.45	1.62	1.52
2	P	2	SER	CB-OG	5.93	1.50	1.42

All (56) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	51	GLU	CA-C-N	-13.22	88.12	117.20
1	D	51	GLU	CA-C-N	-10.92	93.18	117.20
1	D	31	VAL	CG1-CB-CG2	9.57	126.21	110.90
1	D	60	TYR	CB-CG-CD2	-9.52	115.29	121.00
1	D	92	TRP	CD1-CG-CD2	9.36	113.79	106.30
1	B	51	GLU	CA-CB-CG	-9.35	92.82	113.40
1	B	51	GLU	N-CA-C	9.02	135.35	111.00
1	B	54	TYR	CB-CG-CD2	-9.01	115.59	121.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	75	TRP	CD1-CG-CD2	8.78	113.33	106.30
1	B	21	TRP	CD1-CG-CD2	8.77	113.31	106.30
1	B	120	TRP	CD1-CG-CD2	8.43	113.05	106.30
1	B	107	GLN	CA-CB-CG	-8.09	95.59	113.40
1	D	108	TRP	CE2-CD2-CG	-8.04	100.87	107.30
1	D	120	TRP	CD1-CG-CD2	8.01	112.71	106.30
1	D	75	TRP	CD1-CG-CD2	7.96	112.67	106.30
1	D	108	TRP	CD1-CG-CD2	7.89	112.61	106.30
1	B	21	TRP	CE2-CD2-CG	-7.75	101.10	107.30
1	D	79	TRP	CD1-CG-CD2	7.64	112.42	106.30
1	B	108	TRP	CD1-CG-CD2	7.60	112.38	106.30
1	B	120	TRP	CE2-CD2-CG	-7.48	101.32	107.30
1	D	79	TRP	CE2-CD2-CG	-7.25	101.50	107.30
1	B	75	TRP	CE2-CD2-CG	-7.12	101.61	107.30
1	D	54	TYR	CB-CG-CD2	-7.09	116.75	121.00
1	B	108	TRP	CE2-CD2-CG	-7.08	101.64	107.30
1	B	92	TRP	CD1-CG-CD2	7.05	111.94	106.30
1	D	120	TRP	CE2-CD2-CG	-6.97	101.73	107.30
1	D	31	VAL	CB-CA-C	-6.86	98.37	111.40
1	B	133	VAL	CA-CB-CG2	-6.76	100.76	110.90
1	B	92	TRP	CE2-CD2-CG	-6.75	101.90	107.30
1	B	79	TRP	CE2-CD2-CG	-6.74	101.91	107.30
1	D	60	TYR	CB-CG-CD1	6.66	125.00	121.00
1	D	51	GLU	CA-C-O	6.61	133.99	120.10
1	D	21	TRP	CE2-CD2-CG	-6.60	102.02	107.30
1	D	108	TRP	CG-CD2-CE3	6.56	139.80	133.90
1	D	21	TRP	CD1-CG-CD2	6.54	111.53	106.30
1	D	92	TRP	CE2-CD2-CG	-6.45	102.14	107.30
1	B	59	ARG	NE-CZ-NH2	-6.43	117.09	120.30
1	B	79	TRP	CD1-CG-CD2	6.39	111.41	106.30
1	D	46	ALA	O-C-N	-6.33	112.58	122.70
1	D	92	TRP	CG-CD1-NE1	-6.26	103.84	110.10
1	D	75	TRP	CE2-CD2-CG	-6.22	102.33	107.30
1	B	75	TRP	CG-CD1-NE1	-5.99	104.11	110.10
1	D	120	TRP	CG-CD1-NE1	-5.90	104.20	110.10
2	P	6	ASN	O-C-N	5.82	132.01	122.70
1	D	103	ARG	CA-CB-CG	5.81	126.18	113.40
1	D	46	ALA	CA-C-N	5.69	129.72	117.20
1	B	84	ARG	NE-CZ-NH1	5.63	123.12	120.30
1	B	21	TRP	CG-CD2-CE3	5.61	138.95	133.90
1	D	92	TRP	CB-CG-CD1	-5.53	119.81	127.00
1	B	103	ARG	NE-CZ-NH1	5.50	123.05	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	67	ASP	CA-C-N	5.44	127.08	116.20
1	B	21	TRP	CG-CD1-NE1	-5.41	104.69	110.10
2	P	1[A]	PHE	N-CA-C	-5.31	96.67	111.00
2	P	1[B]	PHE	N-CA-C	-5.31	96.67	111.00
1	B	120	TRP	CG-CD1-NE1	-5.20	104.90	110.10
1	D	108	TRP	CB-CG-CD1	-5.06	120.42	127.00

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	134	LYS	Peptide
1	B	51	GLU	Mainchain
1	D	51	GLU	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	B	934	218	878	26	2
1	D	918	219	859	13	0
2	M	48	11	39	2	1
2	P	70	18	61	3	0
3	B	75	150	0	3	1
3	D	80	160	0	0	3
3	M	3	6	0	0	0
3	P	1	2	0	0	0
All	All	2129	784	1837	37	4

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 10.

All (37) 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:22[B]:TYR:CD1	3:B:852:HOH:O	2.19	0.96

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:22[B]:TYR:CE1	1:B:131:THR:HB	2.11	0.86
1:B:114:THR:H	1:D:95:GLN:HE22	1.27	0.81
1:B:51:GLU:HG2	1:B:83:TYR:CD2	2.23	0.73
1:B:22[A]:TYR:HE2	1:B:133:VAL:HG11	1.54	0.72
1:D:13:ALA:N	1:D:96:TYR:HH	1.89	0.69
1:B:44:GLU:HG3	1:B:53:ARG:HG3	1.74	0.68
1:D:51:GLU:HA	1:D:53:ARG:H	1.60	0.65
1:D:120:TRP:CE2	1:D:121:LYS:HD2	2.33	0.64
1:B:51:GLU:HG2	1:B:83:TYR:HD2	1.63	0.62
1:B:95:GLN:HE22	1:D:114:THR:H	1.48	0.62
1:B:22[B]:TYR:CE2	1:B:133:VAL:HB	2.41	0.55
1:B:114:THR:H	1:D:95:GLN:NE2	2.02	0.55
1:B:91:THR:HB	1:D:91:THR:HB	1.93	0.51
1:B:22[A]:TYR:CE2	1:B:133:VAL:HG11	2.41	0.50
1:B:51:GLU:HG3	1:B:81:ASN:ND2	2.27	0.50
1:D:39:LEU:HD13	1:D:60:TYR:CD1	2.48	0.49
1:B:23:ASN:HB3	1:B:130:PHE:CE1	2.48	0.48
1:D:51:GLU:HG3	1:D:81:ASN:ND2	2.28	0.48
1:D:51:GLU:HG2	1:D:83:TYR:CD1	2.49	0.48
1:B:51:GLU:HG3	1:B:81:ASN:HD21	1.79	0.47
1:D:50:ALA:O	1:D:51:GLU:HB2	2.14	0.47
1:B:22[A]:TYR:HE1	1:B:28:THR:HG1	1.61	0.47
1:B:25:LEU:HD12	2:M:6:ASN:CG	2.37	0.45
1:B:133:VAL:HG22	1:B:133:VAL:O	2.16	0.45
1:B:22[B]:TYR:HD1	3:B:852:HOH:O	1.72	0.44
1:B:22[B]:TYR:HE2	1:B:133:VAL:HB	1.79	0.44
1:B:70:GLY:HA3	1:B:95:GLN:HE21	1.83	0.44
1:B:18:THR:HG23	1:B:32:THR:HA	1.99	0.43
1:B:103:ARG:NH1	1:B:131:THR:HG23	2.34	0.42
1:B:13:ALA:N	3:B:984:HOH:O	2.53	0.42
2:P:3:HIS:CE1	2:P:4:PRO:HD2	2.55	0.42
1:B:94:GLY:HA2	1:D:111:THR:HG21	2.02	0.42
1:B:25:LEU:HD12	2:M:6:ASN:ND2	2.34	0.41
1:D:44:GLU:HA	1:D:52:SER:O	2.20	0.41
2:P:1[B]:PHE:HA	2:P:1[B]:PHE:HD1	1.77	0.40

All (4) 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:22[B]:TYR:CE1	3:B:852:HOH:O[3_655]	1.72	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:134:LYS:HZ3	3:D:1062:HOH:H2[8_555]	1.18	0.42
3:D:1056:HOH:O	3:D:1080:HOH:O[4_555]	2.00	0.20
2:M:5:GLN:O	3:D:1025:HOH:H1[6_554]	1.58	0.02

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	B	124/135 (92%)	118 (95%)	5 (4%)	1 (1%)	24	8
1	D	123/135 (91%)	113 (92%)	5 (4%)	5 (4%)	3	0
2	M	4/7 (57%)	4 (100%)	0	0	100	100
2	P	5/7 (71%)	4 (80%)	0	1 (20%)	0	0
All	All	256/284 (90%)	239 (93%)	10 (4%)	7 (3%)	6	0

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	66	THR
1	D	14	GLU
1	D	47	VAL
1	D	51	GLU
1	D	46	ALA
2	P	2	SER
1	D	49	ASN

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	B	93/100 (93%)	84 (90%)	9 (10%)	10	1
1	D	92/100 (92%)	82 (89%)	10 (11%)	8	1
2	M	6/7 (86%)	5 (83%)	1 (17%)	3	0
2	P	8/7 (114%)	8 (100%)	0	100	100
All	All	199/214 (93%)	179 (90%)	20 (10%)	11	1

All (20) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	39	LEU
1	B	45	SER
1	B	49	ASN
1	B	51	GLU
1	B	67	ASP
1	B	69	SER
1	B	73[A]	LEU
1	B	73[B]	LEU
1	B	133	VAL
2	M	2	SER
1	D	31	VAL
1	D	36	ASP
1	D	39	LEU
1	D	45	SER
1	D	51	GLU
1	D	66	THR
1	D	73[A]	LEU
1	D	73[B]	LEU
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 [i](#)

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.