

# Full wwPDB X-ray Structure Validation Report (i)

#### Feb 10, 2024 – 06:53 PM EST

PDB ID : 2RTL

Title: STREPTAVIDIN-2-IMINOBIOTIN-SULFATE COMPLEX, PH 2.50,

SPACE GROUP I4122

Authors : Katz, B.A. Deposited on : 1997-09-11

Resolution : 1.41 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

https://www.wwpdb.org/validation/2017/XrayValidationReportHelp with specific help available everywhere you see the (i) symbol.

The types of validation reports are described at http://www.wwpdb.org/validation/2017/FAQs#types.

The following versions of software and data (see references (1)) were used in the production of this report:

 $Mol Probity \quad : \quad 4.02b\text{--}467$ 

Mogul: 1.8.5 (274361), CSD as541be (2020)

Xtriage (Phenix) : 1.13 EDS : 2.36

Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)

 $Refmac \quad : \quad 5.8.0158$ 

CCP4 : 7.0.044 (Gargrove)

Ideal geometry (proteins) : Engh & Huber (2001) Ideal geometry (DNA, RNA) : Parkinson et al. (1996)

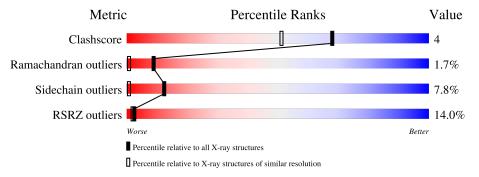
Validation Pipeline (wwPDB-VP) : 2.36

## 1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:  $X\text{-}RAY\ DIFFRACTION$ 

The reported resolution of this entry is 1.41 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	Similar resolution				
Metric	$(\# \mathrm{Entries})$	$(\#  ext{Entries},  ext{ resolution range}(\mathring{ ext{A}}))$				
Clashscore	141614	2696 (1.44-1.40)				
Ramachandran outliers	138981	2632 (1.44-1.40)				
Sidechain outliers	138945	2631 (1.44-1.40)				
RSRZ outliers	127900	2528 (1.44-1.40)				

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 of the lower bar indicate the fraction of residues that contain outliers for >=3, 2, 1 and 0 types of geometric quality criteria respectively. 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 <=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								
			13%								
1	A	135	70%	18%	•	9%					



# 2 Entry composition (i)

There are 4 unique types of molecules in this entry. The entry contains 2174 atoms, of which 1083 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	Λ	193	Total	С	Н	N	О	107	17	0
1	A	123	1989	643	968	171	207	107	11	

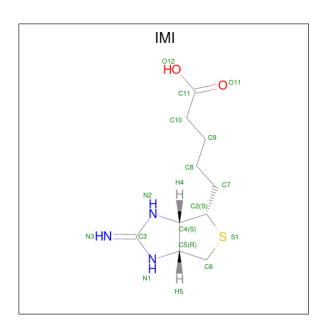
• Molecule 2 is SULFATE ION (three-letter code: SO4) (formula: O<sub>4</sub>S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total 5	O 4	S 1	0	0

 $\bullet$  Molecule 3 is 2-IMINOBIOTIN (three-letter code: IMI) (formula:  $C_{10}H_{17}N_3O_2S).$ 





Mol	Chain	Residues	Atoms				ZeroOcc	AltConf		
2	Λ	1	Total	С	Н	N	О	S	0	0
)	A	A 1	33	10	17	3	2	1	0	0

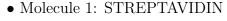
#### • Molecule 4 is water.

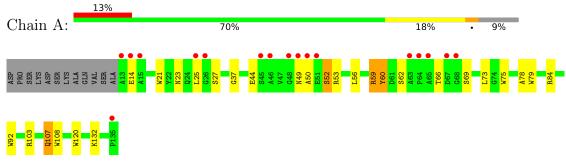
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	49	Total 147	H 98	O 49	0	6



# 3 Residue-property plots (i)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes 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.







# 4 Data and refinement statistics (i)

Property	Value	Source	
Space group	I 41 2 2	Depositor	
Cell constants	58.20Å 58.20Å 173.31Å	Donositor	
a, b, c, $\alpha$ , $\beta$ , $\gamma$	90.00° 90.00° 90.00°	Depositor	
Resolution (Å)	7.50 - 1.41	Depositor	
Resolution (A)	27.59 - 1.30	EDS	
% Data completeness	78.0 (7.50-1.41)	Depositor	
(in resolution range)	67.6 (27.59-1.30)	EDS	
$R_{merge}$	0.06	Depositor	
$R_{sym}$	(Not available)	Depositor	
$< I/\sigma(I) > 1$	1.35 (at 1.30Å)	Xtriage	
Refinement program	X-PLOR	Depositor	
D D.	0.205 , 0.234	Depositor	
$R, R_{free}$	0.209 , (Not available)	DCC	
$R_{free}$ test set	No test flags present.	wwPDB-VP	
Wilson B-factor (Å <sup>2</sup> )	11.3	Xtriage	
Anisotropy	0.249	Xtriage	
Bulk solvent $k_{sol}(e/Å^3)$ , $B_{sol}(Å^2)$	0.23, 65.7	EDS	
L-test for twinning <sup>2</sup>	$ < L > = 0.49, < L^2> = 0.33$	Xtriage	
Estimated twinning fraction	No twinning to report.	Xtriage	
$F_o, F_c$ correlation	0.95	EDS	
Total number of atoms	2174	wwPDB-VP	
Average B, all atoms (Å <sup>2</sup> )	25.0	wwPDB-VP	

Xtriage's analysis on translational NCS is as follows: The largest off-origin peak in the Patterson function is 12.10% of the height of the origin peak. No significant pseudotranslation is detected.

<sup>&</sup>lt;sup>2</sup>Theoretical values of <|L|>,  $<L^2>$  for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.



<sup>&</sup>lt;sup>1</sup>Intensities estimated from amplitudes.

## 5 Model quality (i)

### 5.1 Standard geometry (i)

Bond lengths and bond angles in the following residue types are not validated in this section: SO4, IMI

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	Bo	nd lengths	Bond angles		
		RMSZ	# Z  > 5	RMSZ	# Z  > 5	
1	A	1.33	2/1076~(0.2%)	1.74	30/1471 (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 maintenain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1

#### All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	$\mathbf{Z}$	$Observed(\AA)$	$\operatorname{Ideal}( ext{\AA})$
1	A	120	TRP	CG-CD2	-5.34	1.34	1.43
1	A	21	TRP	CG-CD2	-5.09	1.34	1.43

All (30) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	$\mathbf{Z}$	$Observed(^{o})$	$\mathbf{Ideal}(^{o})$
1	A	108	TRP	CD1-NE1-CE2	9.49	117.54	109.00
1	A	92	TRP	CD1-NE1-CE2	9.45	117.50	109.00
1	A	75	TRP	CD1-NE1-CE2	9.38	117.44	109.00
1	A	120	TRP	CD1-NE1-CE2	9.10	117.19	109.00
1	A	21	TRP	CD1-NE1-CE2	8.88	116.99	109.00
1	A	79	TRP	CD1-NE1-CE2	8.82	116.94	109.00
1	A	120	TRP	NE1-CE2-CZ2	8.56	139.82	130.40
1	A	92	TRP	CG-CD1-NE1	-7.89	102.21	110.10
1	A	108	TRP	NE1-CE2-CZ2	7.78	138.96	130.40
1	A	103	ARG	NE-CZ-NH2	-7.50	116.55	120.30
1	A	21	TRP	NE1-CE2-CZ2	7.46	138.60	130.40

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Mol	Chain	Res	Type	Atoms	$\mathbf{Z}$	$Observed(^o)$	$Ideal(^{o})$
1	A	21	TRP	CG-CD1-NE1	-7.24	102.86	110.10
1	A	120	TRP	CG-CD1-NE1	-7.20	102.90	110.10
1	A	108	TRP	CG-CD1-NE1	-7.12	102.98	110.10
1	A	75	TRP	CG-CD1-NE1	-7.06	103.04	110.10
1	A	53	ARG	NE-CZ-NH2	-6.83	116.89	120.30
1	A	59	ARG	NE-CZ-NH1	6.71	123.66	120.30
1	A	79	TRP	CG-CD1-NE1	-6.44	103.66	110.10
1	A	92	TRP	NE1-CE2-CZ2	6.32	137.35	130.40
1	A	120	TRP	NE1-CE2-CD2	-6.26	101.04	107.30
1	A	108	TRP	NE1-CE2-CD2	-6.07	101.23	107.30
1	A	75	TRP	NE1-CE2-CZ2	5.90	136.89	130.40
1	A	78	ALA	N-CA-C	-5.84	95.22	111.00
1	A	21	TRP	NE1-CE2-CD2	-5.66	101.64	107.30
1	A	92	TRP	NE1-CE2-CD2	-5.51	101.79	107.30
1	A	107[A]	GLN	N-CA-C	-5.47	96.22	111.00
1	A	107[B]	GLN	N-CA-C	-5.47	96.22	111.00
1	A	60[A]	TYR	CB-CG-CD2	-5.43	117.74	121.00
1	A	60[B]	TYR	CB-CG-CD2	-5.43	117.74	121.00
1	A	79	TRP	NE1-CE2-CD2	-5.29	102.01	107.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	84	ARG	Sidechain

### 5.2 Too-close contacts (i)

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	1021	968	950	7	3
2	A	5	0	0	0	0
3	A	16	17	15	2	0
4	A	49	98	0	2	7
All	All	1091	1083	965	7	7

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 (7) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic	Clash
Atom-1	Atom-2	${ m distance}({ m \AA})$	overlap (Å)
1:A:50:ALA:HB2	3:A:300:IMI:H91	1.78	0.65
1:A:52[A]:SER:HB2	4:A:138[A]:HOH:O	2.03	0.58
1:A:132[B]:LYS:NZ	4:A:611:HOH:O	2.39	0.52
1:A:49:ASN:HB3	3:A:300:IMI:C11	2.42	0.49
1:A:37:GLY:HA2	1:A:60[A]:TYR:CE1	2.50	0.47
1:A:23:ASN:ND2	1:A:25:LEU:HD12	2.36	0.40

All (7) 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	$egin{array}{ll}  ext{Interatomic} \  ext{distance } ( ext{Å}) \end{array}$	Clash overlap (Å)
4:A:305:HOH:O	4:A:305:HOH:O[15_555]	0.68	1.52
4:A:305:HOH:O	4:A:305:HOH:H1[15_555]	0.54	1.06
4:A:305:HOH:H1	4:A:305:HOH:H2[15_555]	0.72	0.88
4:A:305:HOH:O	4:A:305:HOH:H2[15_555]	1.04	0.56
1:A:59:ARG:HH21	4:A:231:HOH:H1[15_555]	1.34	0.26
1:A:60[A]:TYR:O	4:A:278:HOH:H1[8_555]	1.53	0.07
1:A:60[B]:TYR:O	4:A:278:HOH:H1[8_555]	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	A	138/135 (102%)	127 (92%)	8 (6%)	3 (2%)	6 0

#### All (3) Ramachandran outliers are listed below:



Mol	Chain	Res	Type
1	A	14	GLU
1	A	52[A]	SER
1	A	52[B]	SER

#### 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	104/100 (104%)	93 (89%)	11 (11%)	6 0

All (11) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	27	SER
1	A	44	GLU
1	A	56	LEU
1	A	62[A]	SER
1	A	62[B]	SER
1	A	66[A]	THR
1	A	66[B]	THR
1	A	69[A]	SER
1	A	69[B]	SER
1	A	107[A]	GLN
1	A	107[B]	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

### 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 monosaccharides in this entry.

### 5.6 Ligand geometry (i)

2 ligands 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).

Mal	Mol Type	Chain	Res	Link	Bo	Bond lengths			Bond angles		
IVIOI	туре	Chain	nes	Lilik	Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2	
2	SO4	A	137	-	4,4,4	0.99	0	6,6,6	0.34	0	
3	IMI	A	300	-	17,17,17	1.52	6 (35%)	20,23,23	2.95	7 (35%)	

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.

$\mathbf{N}$	/Iol	Type	Chain	Res	Link	Chirals	Torsions	Rings
	3	IMI	A	300	-	-	1/7/28/28	0/2/2/2

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(A)	$Ideal(\AA)$
3	A	300	IMI	C2-S1	-2.50	1.78	1.82
3	A	300	IMI	O12-C11	-2.37	1.22	1.30
3	A	300	IMI	C4-N2	2.16	1.49	1.45
3	A	300	IMI	C3-N1	-2.15	1.31	1.35
3	A	300	IMI	C3-N2	-2.15	1.31	1.35
3	A	300	IMI	C5-N1	2.08	1.49	1.46

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	$\mathbf{Z}$	$\operatorname{Observed}(^{o})$	$\operatorname{Ideal}({}^o)$
3	A	300	IMI	C4-C2-S1	6.59	111.49	105.20
3	A	300	IMI	C5-C6-S1	5.65	111.15	106.31

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Mol	Chain	Res	Type	Atoms	$\mathbf{Z}$	$Observed(^o)$	$\operatorname{Ideal}({}^{o})$
3	A	300	IMI	C6-C5-N1	-5.29	106.30	113.03
3	A	300	IMI	C8-C7-C2	-4.79	104.14	113.86
3	A	300	IMI	C2-C4-N2	-4.66	108.95	113.13
3	A	300	IMI	N1-C3-N2	2.58	111.85	109.20
3	A	300	IMI	C2-C4-C5	-2.58	105.94	108.94

There are no chirality outliers.

All (1) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	300	IMI	C4-C2-C7-C8

There are no ring outliers.

1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	300	IMI	2	0

## 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,  $95^{th}$  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	$\langle { m RSRZ} \rangle$	# RSRZ > 2	$OWAB(A^2)$	Q < 0.9
1	A	121/135 (89%)	0.63	17 (14%) 2 2	12, 20, 43, 53	32 (26%)

All (17) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	49	ASN	23.0
1	A	135	PRO	12.5
1	A	50	ALA	8.4
1	A	48	GLY	7.5
1	A	25	LEU	6.8
1	A	46	ALA	6.6
1	A	51	$\operatorname{GLU}$	6.3
1	A	64[A]	PRO	6.1
1	A	65[A]	ALA	5.4
1	A	13	ALA	4.6
1	A	67[A]	ASP	3.6
1	A	68[A]	GLY	3.4
1	A	14	$\operatorname{GLU}$	2.9
1	A	15	ALA	2.7
1	A	63[A]	ALA	2.6
1	A	26	GLY	2.2
1	A	45	SER	2.1

## 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)

There are no monosaccharides in this entry.



# 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. The B-factors column lists the minimum, median,  $95^{th}$  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	$\mathbf{B} ext{-}\mathbf{factors}(\mathbf{\mathring{A}}^2)$	Q<0.9
2	SO4	A	137	5/5	0.92	0.19	40,44,45,48	5
3	IMI	A	300	16/16	0.98	0.05	20,24,39,44	0

### 6.5 Other polymers (i)

There are no such residues in this entry.

