



# Full wwPDB X-ray Structure Validation Report

Feb 27, 2014 – 07:57 PM GMT

PDB ID : 1VWA  
Title : STREPTAVIDIN-FSHPQNT  
Authors : Katz, B.A.; Cass, R.T.  
Deposited on : 1997-03-03  
Resolution : 1.85 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>

---

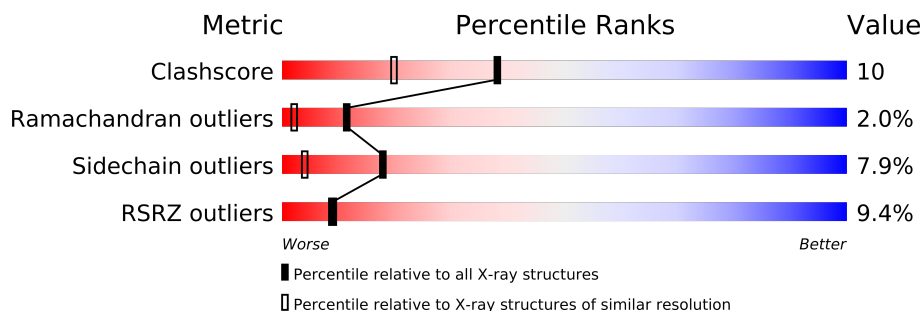
The following versions of software and data (see [references](#)) were used in the production of this report:

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 1.85 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	1470 (1.86-1.86)
Ramachandran outliers	78287	1451 (1.86-1.86)
Sidechain outliers	78261	1451 (1.86-1.86)
RSRZ outliers	66119	1269 (1.86-1.86)

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	
1	D	123	
2	M	7	
2	P	7	

## 2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 2957 atoms, of which 824 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	0	1	0
			1146	580	220	159	187			
1	D	121	Total	C	H	N	O	14	2	0
			1127	569	217	157	184			

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

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

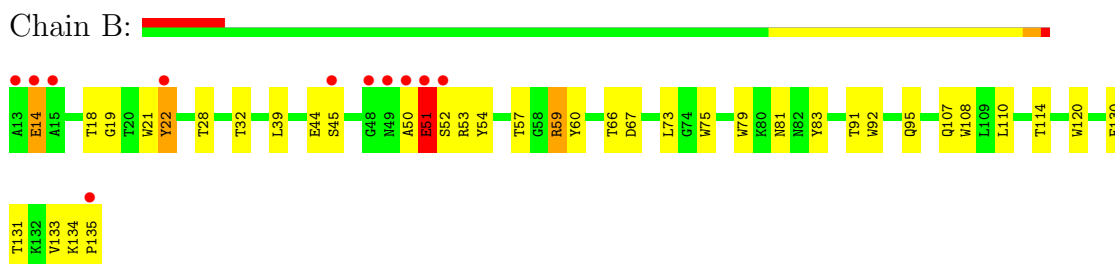
- Molecule 3 is water.

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	B	89	Total	H	O	0	0
			267	178	89		
3	D	85	Total	H	O	0	0
			255	170	85		
3	M	3	Total	H	O	0	0
			9	6	3		
3	P	2	Total	H	O	0	0
			6	4	2		

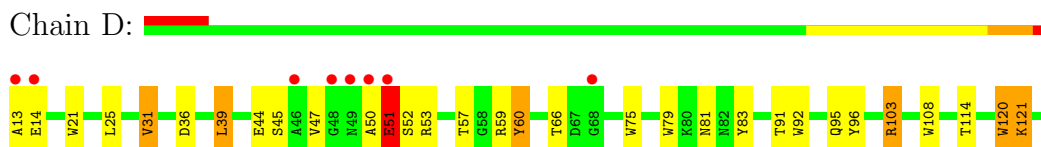
### 3 Residue-property plots

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



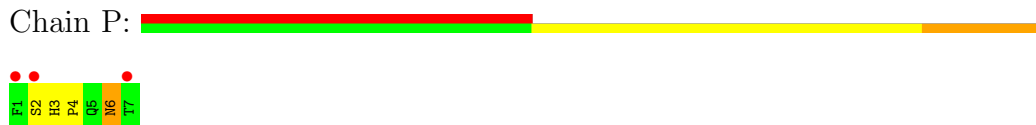
- Molecule 1: STREPTAVIDIN



- Molecule 2: PEPTIDE LIGAND CONTAINING HPQ



- Molecule 2: PEPTIDE LIGAND CONTAINING HPQ



## 4 Data and refinement statistics

Property	Value	Source
Space group	I 2 2 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	97.03Å 107.47Å 48.55Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	7.50 – 1.85 30.97 – 1.80	Depositor EDS
% Data completeness (in resolution range)	71.4 (7.50-1.85) 69.4 (30.97-1.80)	Depositor EDS
$R_{merge}$	0.08	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.94 (at 1.80Å)	Xtriage
Refinement program	X-PLOR	Depositor
R, $R_{free}$	0.192 , 0.259 0.197 , (Not available)	Depositor DCC
$R_{free}$ test set	No test flags present.	DCC
Wilson B-factor (Å <sup>2</sup> )	13.6	Xtriage
Anisotropy	0.057	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.34 , 64.9	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning	$\langle  L  \rangle = 0.48$ , $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 16643 reflections	Xtriage
$F_o, F_c$ correlation	0.92	EDS
Total number of atoms	2957	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	17.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.54% 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

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	0/954	1.80	28/1305 (2.1%)
1	D	0.89	0/940	1.77	30/1287 (2.3%)
2	M	0.87	0/49	1.39	0/65
2	P	1.01	0/73	1.89	1/97 (1.0%)
All	All	0.93	0/2016	1.78	59/2754 (2.1%)

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

There are no bond length outliers.

All (59) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	51	GLU	CA-C-N	-12.68	89.31	117.20
1	B	75	TRP	CD1-CG-CD2	9.12	113.60	106.30
1	D	92	TRP	CD1-CG-CD2	9.13	113.60	106.30
1	B	120	TRP	CD1-CG-CD2	8.81	113.35	106.30
1	D	75	TRP	CD1-CG-CD2	8.50	113.10	106.30
1	B	51	GLU	CA-CB-CG	-8.46	94.78	113.40
1	B	75	TRP	CE2-CD2-CG	-8.34	100.63	107.30
1	D	108	TRP	CD1-CG-CD2	8.21	112.86	106.30
1	D	120	TRP	CD1-CG-CD2	8.20	112.86	106.30
1	B	108	TRP	CD1-CG-CD2	8.02	112.72	106.30
1	B	21	TRP	CD1-CG-CD2	7.96	112.67	106.30
1	B	108	TRP	CE2-CD2-CG	-7.88	101.00	107.30
1	B	92	TRP	CD1-CG-CD2	7.84	112.58	106.30
1	D	108	TRP	CE2-CD2-CG	-7.83	101.04	107.30

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	79	TRP	CE2-CD2-CG	-7.80	101.06	107.30
1	D	75	TRP	CE2-CD2-CG	-7.70	101.14	107.30
1	B	51	GLU	N-CA-C	7.61	131.56	111.00
1	B	92	TRP	CE2-CD2-CG	-7.57	101.24	107.30
1	B	120	TRP	CE2-CD2-CG	-7.43	101.35	107.30
1	D	79	TRP	CD1-CG-CD2	7.43	112.24	106.30
1	B	21	TRP	CE2-CD2-CG	-7.18	101.55	107.30
1	D	21	TRP	CE2-CD2-CG	-7.17	101.56	107.30
1	D	21	TRP	CD1-CG-CD2	7.13	112.00	106.30
1	D	51	GLU	CA-C-N	-7.05	101.68	117.20
1	D	92	TRP	CE2-CD2-CG	-6.99	101.71	107.30
1	D	120	TRP	CE2-CD2-CG	-6.93	101.75	107.30
1	B	79	TRP	CD1-CG-CD2	6.84	111.77	106.30
1	D	31	VAL	CB-CA-C	-6.79	98.50	111.40
1	D	31	VAL	CG1-CB-CG2	6.61	121.48	110.90
1	D	92	TRP	CB-CG-CD1	-6.61	118.41	127.00
1	D	60	TYR	CB-CG-CD2	-6.53	117.08	121.00
1	B	79	TRP	CE2-CD2-CG	-6.39	102.19	107.30
2	P	6	ASN	O-C-N	6.19	132.60	122.70
1	B	108	TRP	CG-CD2-CE3	6.10	139.39	133.90
1	D	92	TRP	CG-CD1-NE1	-6.03	104.07	110.10
1	D	120	TRP	CG-CD1-NE1	-5.95	104.15	110.10
1	D	57	THR	CA-CB-CG2	5.95	120.72	112.40
1	B	75	TRP	CG-CD1-NE1	-5.94	104.16	110.10
1	B	75	TRP	CG-CD2-CE3	5.83	139.15	133.90
1	B	22[A]	TYR	CB-CG-CD1	-5.82	117.51	121.00
1	B	22[B]	TYR	CB-CG-CD1	-5.82	117.51	121.00
1	D	75	TRP	CG-CD2-CE3	5.80	139.12	133.90
1	D	108	TRP	CG-CD2-CE3	5.72	139.04	133.90
1	D	21	TRP	CG-CD2-CE3	5.71	139.04	133.90
1	D	75	TRP	CB-CG-CD1	-5.69	119.60	127.00
1	B	120	TRP	CG-CD1-NE1	-5.67	104.43	110.10
1	D	103	ARG	CA-CB-CG	5.66	125.84	113.40
1	D	79	TRP	CG-CD2-CE3	5.56	138.91	133.90
1	B	108	TRP	CB-CG-CD1	-5.50	119.86	127.00
1	B	54	TYR	CB-CG-CD2	-5.45	117.73	121.00
1	D	92	TRP	CG-CD2-CE3	5.40	138.76	133.90
1	B	107	GLN	CA-CB-CG	-5.37	101.60	113.40
1	B	59	ARG	NE-CZ-NH2	-5.35	117.63	120.30
1	B	39	LEU	CA-CB-CG	5.34	127.58	115.30
1	D	21	TRP	CB-CG-CD1	-5.09	120.38	127.00
1	D	108	TRP	CB-CG-CD1	-5.09	120.38	127.00

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	51	GLU	CA-C-O	5.04	130.69	120.10
1	B	21	TRP	CG-CD1-NE1	-5.03	105.07	110.10
1	B	92	TRP	CG-CD2-CE3	5.01	138.41	133.90

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	134	LYS	Peptide
1	B	51	GLU	Mainchain

## 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	926	220	648	25	5
1	D	910	217	638	17	0
2	M	48	11	28	0	0
2	P	70	18	44	2	0
3	B	89	178	0	4	1
3	D	85	170	0	0	5
3	M	3	6	0	0	0
3	P	2	4	0	0	0
All	All	2133	824	1358	37	6

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 10.

All (37) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:22[B]:TYR:CD1	3:B:852:HOH:O	2.25	0.87
1:B:22[B]:TYR:CE1	1:B:131:THR:HB	2.15	0.80
1:B:59:ARG:HH11	3:B:643:HOH:H1	1.31	0.76
1:B:51:GLU:HG2	1:B:83:TYR:CD1	2.25	0.70
3:B:605:HOH:H1	1:D:59:ARG:HE	1.39	0.69

*Continued on next page...*



Continued from previous page...

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:22[B]:TYR:HD1	3:B:852:HOH:O	1.71	0.68
1:B:22[B]:TYR:HE2	1:B:133:VAL:HB	1.57	0.68
1:B:22[A]:TYR:HE2	1:B:133:VAL:HG11	1.60	0.67
1:B:22[B]:TYR:CE2	1:B:133:VAL:HB	2.32	0.65
1:B:14:GLU:HB3	1:B:60:TYR:OH	2.02	0.58
1:B:51:GLU:HG2	1:B:83:TYR:HD1	1.64	0.58
1:D:44:GLU:HB3	1:D:53:ARG:HG2	1.86	0.58
1:B:114:THR:H	1:D:95:GLN:NE2	2.03	0.56
1:D:51:GLU:HG3	1:D:81:ASN:ND2	2.23	0.53
1:B:51:GLU:HG3	1:B:81:ASN:ND2	2.23	0.53
1:D:51:GLU:HG2	1:D:83:TYR:CD2	2.45	0.52
1:B:50:ALA:O	1:B:53:ARG:HB2	2.09	0.52
1:B:114:THR:H	1:D:95:GLN:HE21	1.59	0.51
1:D:51:GLU:HG2	1:D:83:TYR:CE2	2.46	0.50
1:B:50:ALA:HA	1:B:53:ARG:HH21	1.77	0.50
1:D:44:GLU:HA	1:D:52:SER:O	2.13	0.48
1:D:120:TRP:CE2	1:D:121:LYS:HD2	2.48	0.48
1:D:13:ALA:N	1:D:96:TYR:HH	2.13	0.47
1:D:13:ALA:H3	1:D:96:TYR:HE2	1.64	0.46
1:B:95:GLN:NE2	1:D:114:THR:H	2.14	0.45
1:D:25:LEU:HD12	2:P:6:ASN:CG	2.37	0.45
1:B:91:THR:HB	1:D:91:THR:HB	1.98	0.44
1:B:22[A]:TYR:CE2	1:B:133:VAL:HG11	2.47	0.44
1:B:44:GLU:HA	1:B:52:SER:O	2.18	0.43
2:P:3:HIS:CE1	2:P:4:PRO:HD2	2.53	0.43
1:B:95:GLN:HE21	1:D:114:THR:H	1.66	0.43
1:D:39:LEU:HD13	1:D:60:TYR:CD1	2.53	0.42
1:B:18:THR:HG23	1:B:32:THR:HA	1.99	0.42
1:B:19:GLY:HA2	1:B:135:PRO:HA	2.01	0.42
1:B:22[A]:TYR:HE1	1:B:28:THR:OG1	2.04	0.41
1:B:57:THR:HG22	1:D:59:ARG:HG2	2.02	0.41

All (6) 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	Distance(Å)	Clash(Å)
3:D:585:HOH:O	3:D:585:HOH:O[3_556]	1.54	0.66
1:B:22[A]:TYR:OH	3:D:585:HOH:O[8_555]	1.76	0.44
1:B:22[B]:TYR:CE1	3:B:852:HOH:O[3_655]	1.78	0.42
1:B:22[A]:TYR:CE2	3:D:1095:HOH:O[6_554]	1.92	0.28
1:B:22[A]:TYR:CD2	3:D:1095:HOH:O[6_554]	1.93	0.27
1:B:22[A]:TYR:CZ	3:D:585:HOH:O[8_555]	2.19	0.01

## 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	122/123 (99%)	116 (95%)	5 (4%)	1 (1%)	27	10
1	D	121/123 (98%)	112 (93%)	6 (5%)	3 (2%)	9	1
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	252/260 (97%)	236 (94%)	11 (4%)	5 (2%)	11	2

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	66	THR
1	D	14	GLU
1	D	47	VAL
1	D	50	ALA
2	P	2	SER

### 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	91/90 (101%)	85 (93%)	6 (7%)	24	7
1	D	90/90 (100%)	82 (91%)	8 (9%)	14	2
2	M	6/7 (86%)	5 (83%)	1 (17%)	3	0
2	P	8/7 (114%)	8 (100%)	0	100	100
All	All	195/194 (100%)	180 (92%)	15 (8%)	18	4

All (15) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	14	GLU
1	B	45	SER
1	B	51	GLU
1	B	67	ASP
1	B	73	LEU
1	B	110	LEU
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	103	ARG
1	D	121	LYS
2	M	2	SER

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 ⓘ

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	123/123 (100%)	0.30	11 (8%)	10 10	5, 17, 49, 66	14 (11%)
1	D	119/123 (96%)	0.36	8 (6%)	17 16	5, 15, 45, 55	9 (7%)
2	M	5/7 (71%)	3.50	2 (40%)	1 1	34, 43, 52, 54	1 (20%)
2	P	7/7 (100%)	1.93	3 (42%)	1 1	34, 44, 68, 72	0
All	All	254/260 (97%)	0.43	24 (9%)	9 9	5, 16, 50, 72	24 (9%)

All (24) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	14	GLU	10.4
2	M	2	SER	10.0
1	D	46	ALA	8.6
1	B	48	GLY	8.5
1	D	48	GLY	7.5
1	B	14	GLU	7.4
1	B	49	ASN	4.8
1	B	15	ALA	4.7
2	P	2	SER	4.7
1	D	50	ALA	4.6
1	B	13	ALA	4.1
2	M	4	PRO	4.1
1	D	13	ALA	4.0
1	B	22[A]	TYR	3.9
1	B	50	ALA	3.5
1	D	51	GLU	3.5
2	P	1[A]	PHE	3.2
1	B	135	PRO	3.1
2	P	7	THR	3.0
1	D	68	GLY	2.8
1	B	52	SER	2.8

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Res	Type	RSRZ
1	B	45	SER	2.4
1	B	51	GLU	2.4
1	D	49	ASN	2.0

## 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.