



# Full wwPDB X-ray Structure Validation Report

Feb 26, 2014 – 11:57 PM GMT

PDB ID : 1SH5  
Title : Crystal structure of actin-binding domain of mouse plectin  
Authors : Sevcik, J.; Urbanikova, L.  
Deposited on : 2004-02-25  
Resolution : 2.00 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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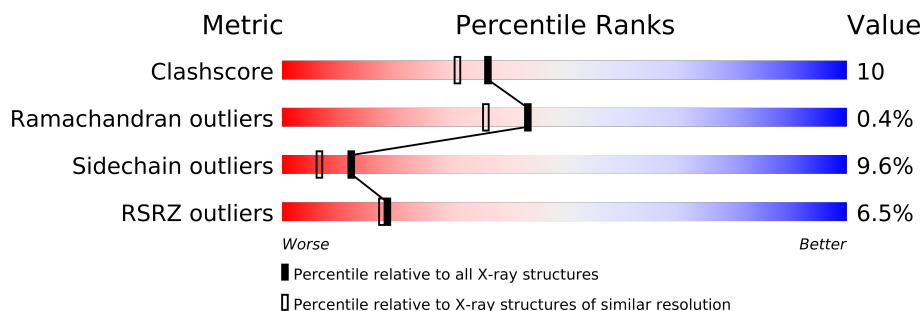
The following versions of software and data (see [references](#)) were used in the production of this report:

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 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	79885	6188 (2.00-2.00)
Ramachandran outliers	78287	6102 (2.00-2.00)
Sidechain outliers	78261	6100 (2.00-2.00)
RSRZ outliers	66119	4890 (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  $\geq 3$ , 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	A	245	
1	B	245	

## 2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 8252 atoms, of which 4226 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 Plectin 1.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	232	Total	C	H	N	O	S	138	0	0
			3844	1209	1925	350	353	7			
1	B	232	Total	C	H	N	O	S	138	0	0
			3844	1209	1925	350	353	7			

There are 16 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	GLY	-	CLONING ARTIFACT	UNP Q9QXS1
A	2	SER	-	CLONING ARTIFACT	UNP Q9QXS1
A	3	HIS	-	CLONING ARTIFACT	UNP Q9QXS1
A	4	MET	-	CLONING ARTIFACT	UNP Q9QXS1
A	5	GLU	-	CLONING ARTIFACT	UNP Q9QXS1
A	6	PHE	-	CLONING ARTIFACT	UNP Q9QXS1
A	244	GLU	-	CLONING ARTIFACT	UNP Q9QXS1
A	245	PHE	-	CLONING ARTIFACT	UNP Q9QXS1
B	1	GLY	-	CLONING ARTIFACT	UNP Q9QXS1
B	2	SER	-	CLONING ARTIFACT	UNP Q9QXS1
B	3	HIS	-	CLONING ARTIFACT	UNP Q9QXS1
B	4	MET	-	CLONING ARTIFACT	UNP Q9QXS1
B	5	GLU	-	CLONING ARTIFACT	UNP Q9QXS1
B	6	PHE	-	CLONING ARTIFACT	UNP Q9QXS1
B	244	GLU	-	CLONING ARTIFACT	UNP Q9QXS1
B	245	PHE	-	CLONING ARTIFACT	UNP Q9QXS1

- Molecule 2 is water.

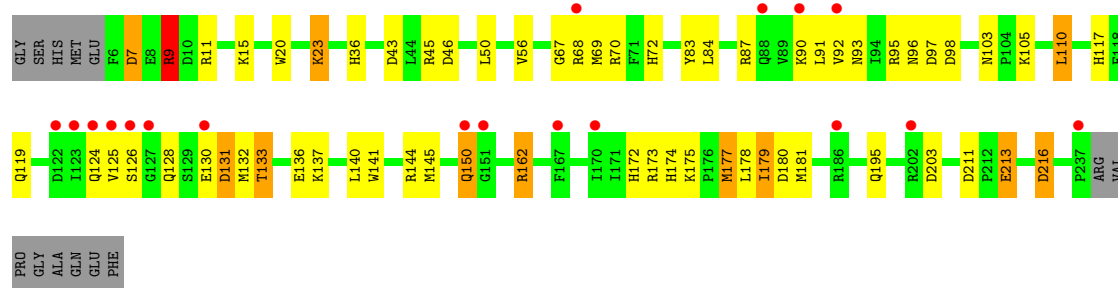
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	96	Total	H	O	192	0
			288	192	96		
2	B	92	Total	H	O	184	0
			276	184	92		

### 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.

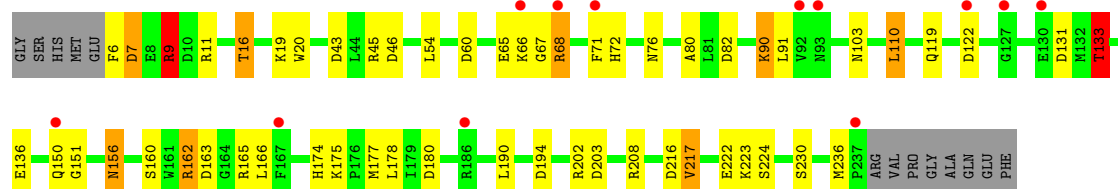
#### • Molecule 1: Plectin 1

Chain A:



#### • Molecule 1: Plectin 1

Chain B:



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	55.31Å 108.92Å 63.75Å 90.00° 115.25° 90.00°	Depositor
Resolution (Å)	30.00 – 2.00 19.97 – 1.95	Depositor EDS
% Data completeness (in resolution range)	97.9 (30.00-2.00) 91.8 (19.97-1.95)	Depositor EDS
$R_{merge}$	0.06	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.71 (at 1.94Å)	Xtriage
Refinement program	REFMAC 5.1.24	Depositor
R, $R_{free}$	0.151 , 0.194 0.163 , (Not available)	Depositor DCC
$R_{free}$ test set	No test flags present.	DCC
Wilson B-factor (Å <sup>2</sup> )	31.0	Xtriage
Anisotropy	0.495	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.38 , 43.7	EDS
Estimated twinning fraction	0.028 for h,-k,-h-l	Xtriage
L-test for twinning	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 47777 reflections	Xtriage
$F_o, F_c$ correlation	0.97	EDS
Total number of atoms	8252	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	43.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.24% 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	A	1.17	2/1957 (0.1%)	1.25	17/2643 (0.6%)
1	B	1.22	9/1957 (0.5%)	1.33	24/2643 (0.9%)
All	All	1.19	11/3914 (0.3%)	1.29	41/5286 (0.8%)

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	1

All (11) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	9	ARG	CD-NE	-9.77	1.29	1.46
1	A	9	ARG	CD-NE	-7.49	1.33	1.46
1	B	9	ARG	CZ-NH2	-6.67	1.24	1.33
1	B	177	MET	CG-SD	6.54	1.98	1.81
1	B	222	GLU	CD-OE1	-6.29	1.18	1.25
1	A	56	VAL	CB-CG1	-6.07	1.40	1.52
1	B	80	ALA	CA-CB	5.33	1.63	1.52
1	B	230	SER	CB-OG	-5.21	1.35	1.42
1	B	19	LYS	CE-NZ	-5.11	1.36	1.49
1	B	223	LYS	CD-CE	5.06	1.63	1.51
1	B	222	GLU	CD-OE2	-5.02	1.20	1.25

All (41) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	9	ARG	NE-CZ-NH2	-25.05	107.77	120.30
1	A	9	ARG	NE-CZ-NH2	-21.70	109.45	120.30
1	B	9	ARG	NE-CZ-NH1	19.02	129.81	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	9	ARG	NE-CZ-NH1	14.50	127.55	120.30
1	B	203	ASP	CB-CG-OD2	10.88	128.09	118.30
1	B	11	ARG	NE-CZ-NH1	-10.37	115.12	120.30
1	A	45	ARG	NE-CZ-NH1	-8.51	116.05	120.30
1	A	203	ASP	CB-CG-OD2	8.31	125.78	118.30
1	B	163	ASP	CB-CG-OD2	8.23	125.70	118.30
1	B	46	ASP	CB-CG-OD1	8.15	125.64	118.30
1	A	46	ASP	CB-CG-OD1	8.01	125.51	118.30
1	A	98	ASP	CB-CG-OD1	7.47	125.02	118.30
1	A	11	ARG	NE-CZ-NH1	-7.45	116.58	120.30
1	B	180	ASP	CB-CG-OD2	7.29	124.86	118.30
1	A	7	ASP	CB-CG-OD1	7.22	124.80	118.30
1	B	11	ARG	NE-CZ-NH2	7.14	123.87	120.30
1	B	45	ARG	NE-CZ-NH2	6.55	123.58	120.30
1	B	60	ASP	CB-CG-OD2	6.49	124.14	118.30
1	B	203	ASP	CB-CG-OD1	-6.47	112.47	118.30
1	B	16	THR	CA-CB-OG1	6.36	122.36	109.00
1	B	133	THR	N-CA-CB	-6.22	98.49	110.30
1	B	7	ASP	CB-CG-OD1	6.18	123.86	118.30
1	A	131	ASP	CB-CG-OD2	6.16	123.84	118.30
1	A	97	ASP	CB-CG-OD2	6.13	123.81	118.30
1	B	82	ASP	CB-CG-OD2	5.84	123.56	118.30
1	B	43	ASP	CB-CG-OD2	5.83	123.55	118.30
1	B	217	VAL	CG1-CB-CG2	5.80	120.17	110.90
1	B	165	ARG	NE-CZ-NH2	-5.74	117.43	120.30
1	A	11	ARG	NE-CZ-NH2	5.71	123.15	120.30
1	A	180	ASP	CB-CG-OD2	5.70	123.43	118.30
1	A	43	ASP	CB-CG-OD2	5.64	123.37	118.30
1	B	194	ASP	CB-CG-OD1	5.55	123.30	118.30
1	A	110	LEU	CB-CG-CD2	5.52	120.38	111.00
1	B	223	LYS	CD-CE-NZ	-5.46	99.14	111.70
1	A	216	ASP	CB-CG-OD2	5.26	123.04	118.30
1	B	122	ASP	CB-CG-OD2	5.23	123.01	118.30
1	B	131	ASP	CB-CG-OD2	5.15	122.94	118.30
1	A	177	MET	CG-SD-CE	5.13	108.41	100.20
1	B	45	ARG	NE-CZ-NH1	-5.08	117.76	120.30
1	A	45	ARG	NE-CZ-NH2	5.02	122.81	120.30
1	B	110	LEU	CB-CG-CD2	5.01	119.52	111.00

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	151	GLY	Peptide

## 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	A	1919	1925	0	42	0
1	B	1919	1925	0	35	0
2	A	96	192	0	12	0
2	B	92	184	0	8	0
All	All	4026	4226	0	76	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 10.

All (76) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:69:MET:CE	1:A:69:MET:SD	2.03	1.47
1:B:16:THR:HG23	1:B:16:THR:OG1	0.95	1.14
1:B:16:THR:OG1	1:B:16:THR:HG22	0.96	1.13
1:B:16:THR:OG1	1:B:16:THR:HG21	0.95	1.12
1:B:16:THR:HG22	1:B:16:THR:CB	1.89	1.02
1:B:16:THR:HG21	1:B:16:THR:CB	1.90	1.02
1:B:16:THR:HG23	1:B:16:THR:CB	1.90	1.01
1:B:162:ARG:NH1	1:B:216:ASP:OD1	2.05	0.89
1:B:16:THR:HG21	1:B:16:THR:HG22	1.55	0.88
1:B:16:THR:HG23	1:B:16:THR:HG21	1.55	0.88
1:B:16:THR:HG23	1:B:16:THR:HG22	1.56	0.88
1:A:213:GLU:HG3	2:A:329:HOH:O	1.76	0.86
1:A:23:LYS:CD	2:A:339:HOH:O	2.28	0.81
1:A:23:LYS:HD3	2:A:339:HOH:O	1.80	0.80
1:A:15:LYS:NZ	1:A:211:ASP:OD2	2.14	0.79
1:A:23:LYS:HE3	2:A:339:HOH:O	1.83	0.77
1:A:162:ARG:NH2	1:A:216:ASP:OD1	2.19	0.75
1:A:23:LYS:CE	2:A:339:HOH:O	2.39	0.70
1:A:124:GLN:HA	1:A:128:GLN:NE2	2.06	0.69

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:7:ASP:OD1	1:A:9:ARG:HD3	1.93	0.68
1:A:7:ASP:OD1	1:A:9:ARG:CD	2.42	0.68
1:B:202:ARG:HD2	2:B:297:HOH:O	1.97	0.63
1:B:16:THR:HG22	1:B:16:THR:CG2	2.29	0.63
1:A:125:VAL:O	1:A:128:GLN:HG3	1.99	0.62
1:B:65:GLU:H	1:B:76:ASN:ND2	1.97	0.61
1:B:20:TRP:HE1	1:B:119:GLN:NE2	1.97	0.61
1:A:162:ARG:HH21	1:A:162:ARG:CG	2.13	0.61
1:B:67:GLY:HA3	1:B:72:HIS:CD2	2.36	0.61
1:A:179:ILE:HD12	1:A:181:MET:SD	2.42	0.60
1:A:126:SER:O	1:A:144:ARG:NH2	2.30	0.59
1:A:145:MET:CE	1:A:174:HIS:CE1	2.86	0.58
1:B:16:THR:HG23	1:B:16:THR:HB	1.84	0.58
1:A:20:TRP:HE1	1:A:119:GLN:NE2	2.02	0.57
1:A:162:ARG:HH21	1:A:162:ARG:HG2	1.70	0.56
1:B:16:THR:HG21	2:B:252:HOH:O	2.05	0.56
1:B:133:THR:HG21	1:B:136:GLU:H	1.72	0.55
1:B:174:HIS:CD2	1:B:236:MET:HG2	2.42	0.55
1:B:175:LYS:HE3	2:B:282:HOH:O	2.07	0.54
1:B:71:PHE:CB	2:B:288:HOH:O	2.55	0.53
1:A:213:GLU:CG	2:A:329:HOH:O	2.43	0.53
1:B:156:ASN:ND2	1:B:160:SER:H	2.07	0.53
1:A:145:MET:HE2	1:A:174:HIS:CE1	2.44	0.52
1:B:162:ARG:CG	1:B:162:ARG:HH11	2.22	0.52
1:A:69:MET:H	1:A:72:HIS:HD2	1.57	0.52
1:A:133:THR:HG21	1:A:136:GLU:H	1.75	0.51
1:B:7:ASP:OD1	1:B:9:ARG:CD	2.58	0.51
1:A:36:HIS:HD2	2:A:307:HOH:O	1.94	0.51
1:A:213:GLU:CD	1:A:213:GLU:H	2.14	0.51
1:A:141:TRP:O	1:A:145:MET:HG2	2.11	0.50
1:A:211:ASP:HB3	1:A:213:GLU:OE1	2.11	0.50
1:A:9:ARG:HD2	2:A:331:HOH:O	2.10	0.50
1:B:16:THR:HG23	1:B:224:SER:OG	2.12	0.49
1:B:68:ARG:N	1:B:72:HIS:HD2	2.11	0.49
1:A:67:GLY:HA3	1:A:72:HIS:CD2	2.47	0.49
1:A:213:GLU:HG3	2:A:306:HOH:O	2.13	0.48
1:A:131:ASP:HB2	1:B:90:LYS:NZ	2.28	0.48
1:B:67:GLY:C	1:B:72:HIS:CD2	2.87	0.48
1:B:174:HIS:HE1	2:B:282:HOH:O	1.95	0.47
1:B:71:PHE:HB2	2:B:288:HOH:O	2.15	0.47
1:A:83:TYR:CZ	1:A:87:ARG:HD2	2.49	0.47
1:A:145:MET:HE1	1:A:174:HIS:CE1	2.50	0.47

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:179:ILE:CD1	1:A:181:MET:SD	3.03	0.46
1:A:172:HIS:CD2	1:A:181:MET:HG2	2.50	0.46
1:A:23:LYS:HG2	2:A:295:HOH:O	2.16	0.45
2:A:290:HOH:O	1:B:133:THR:HG22	2.18	0.44
1:B:67:GLY:CA	1:B:72:HIS:CD2	3.01	0.44
1:A:132:MET:CE	1:A:137:LYS:HG2	2.48	0.43
1:A:87:ARG:NH1	1:A:117:HIS:NE2	2.66	0.43
1:B:6:PHE:N	2:B:323:HOH:O	2.51	0.43
1:A:92:VAL:O	1:A:93:ASN:HB2	2.19	0.42
1:A:145:MET:CE	1:A:174:HIS:HE1	2.33	0.42
1:B:71:PHE:HB3	2:B:288:HOH:O	2.20	0.41
1:B:7:ASP:OD1	1:B:9:ARG:HD3	2.20	0.41
1:A:36:HIS:HE1	2:A:318:HOH:O	2.03	0.41
1:A:83:TYR:CE2	1:A:87:ARG:HD2	2.55	0.41
1:A:172:HIS:HB2	1:A:179:ILE:CG1	2.52	0.40

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	A	230/245 (94%)	225 (98%)	3 (1%)	2 (1%)	25	14
1	B	230/245 (94%)	228 (99%)	2 (1%)	0	100	100
All	All	460/490 (94%)	453 (98%)	5 (1%)	2 (0%)	43	36

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	150	GLN
1	A	96	ASN

### 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	A	213/223 (96%)	189 (89%)	24 (11%)	9	4
1	B	213/223 (96%)	196 (92%)	17 (8%)	17	10
All	All	426/446 (96%)	385 (90%)	41 (10%)	12	6

All (41) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	9	ARG
1	A	23	LYS
1	A	50	LEU
1	A	68	ARG
1	A	70	ARG
1	A	84	LEU
1	A	90	LYS
1	A	91	LEU
1	A	95	ARG
1	A	103	ASN
1	A	105	LYS
1	A	110	LEU
1	A	130	GLU
1	A	133	THR
1	A	140	LEU
1	A	150	GLN
1	A	162	ARG
1	A	173	ARG
1	A	175	LYS
1	A	177	MET
1	A	178	LEU
1	A	179	ILE
1	A	195	GLN
1	A	213	GLU
1	B	9	ARG
1	B	54	LEU
1	B	66	LYS
1	B	68	ARG
1	B	90	LYS

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Mol	Chain	Res	Type
1	B	91	LEU
1	B	103	ASN
1	B	110	LEU
1	B	133	THR
1	B	150	GLN
1	B	156	ASN
1	B	162	ARG
1	B	166	LEU
1	B	178	LEU
1	B	190	LEU
1	B	208	ARG
1	B	217	VAL

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (11) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	22	ASN
1	A	72	HIS
1	A	76	ASN
1	A	96	ASN
1	A	128	GLN
1	A	156	ASN
1	A	174	HIS
1	B	22	ASN
1	B	72	HIS
1	B	78	GLN
1	B	96	ASN

### 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	A	232/245 (94%)	0.16	18 (7%) 13 12	19, 42, 72, 92	0
1	B	232/245 (94%)	0.08	12 (5%) 26 26	20, 39, 69, 78	0
All	All	464/490 (94%)	0.12	30 (6%) 18 18	19, 40, 71, 92	0

All (30) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	92	VAL	6.2
1	A	92	VAL	4.5
1	B	237	PRO	4.4
1	B	71	PHE	4.1
1	A	237	PRO	3.9
1	A	127	GLY	3.6
1	B	130	GLU	3.4
1	A	186	ARG	3.4
1	A	122	ASP	3.3
1	A	150	GLN	3.3
1	A	68	ARG	3.1
1	B	68	ARG	2.8
1	A	167	PHE	2.8
1	B	66	LYS	2.8
1	B	93	ASN	2.8
1	B	150	GLN	2.7
1	A	202	ARG	2.7
1	A	123	ILE	2.7
1	A	126	SER	2.6
1	A	151	GLY	2.5
1	B	167	PHE	2.5
1	B	122	ASP	2.5
1	A	130	GLU	2.4
1	A	90	LYS	2.4

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Mol	Chain	Res	Type	RSRZ
1	A	125	VAL	2.4
1	A	88	GLN	2.2
1	A	124	GLN	2.2
1	B	127	GLY	2.1
1	A	170	ILE	2.1
1	B	186	ARG	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.