



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 06:53 AM UTC

PDB ID : 3ZU7 / pdb_00003zu7
Title : Crystal structure of a designed selected Ankyrin Repeat protein in complex with the MAP kinase ERK2
Authors : Kummer, L.; Mittl, P.R.; Pluckthun, A.
Deposited on : 2011-07-16
Resolution : 1.97 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 ⓘ 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 ⓘ](#)) were used in the production of this report:

MolProbity	:	4-5-2 with Phenix2.0
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

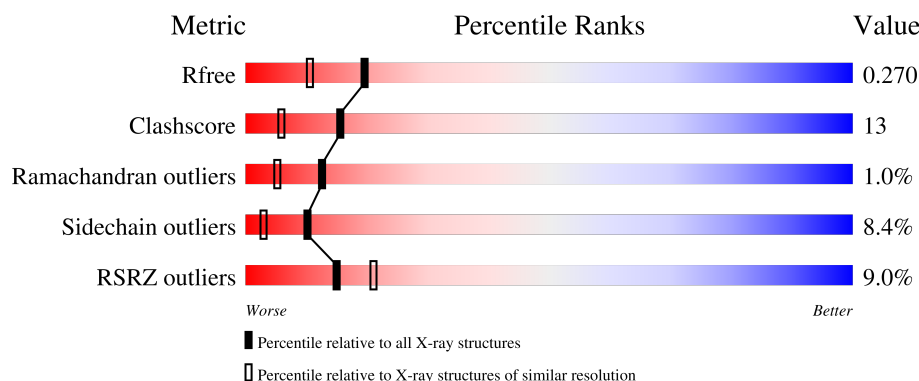
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.97 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))
R_{free}	180053	1506 (1.98-1.98)
Clashscore	190562	1534 (1.98-1.98)
Ramachandran outliers	187476	1518 (1.98-1.98)
Sidechain outliers	187428	1518 (1.98-1.98)
RSRZ outliers	180081	1506 (1.98-1.98)

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 $\leq 5\%$. 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.

Mol	Chain	Length	Quality of chain
1	A	365	10% 65% 22% 6% 6%
2	B	169	4% 78% 11% 9%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 4220 atoms, of which 0 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 MITOGEN-ACTIVATED PROTEIN KINASE 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	344	Total	C	N	O	S	0	0	0
			2817	1809	481	512	15			

There are 9 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-6	ALA	-	expression tag	UNP P63086
A	-5	HIS	-	expression tag	UNP P63086
A	-4	HIS	-	expression tag	UNP P63086
A	-3	HIS	-	expression tag	UNP P63086
A	-2	HIS	-	expression tag	UNP P63086
A	-1	HIS	-	expression tag	UNP P63086
A	0	HIS	-	expression tag	UNP P63086
A	1	ALA	-	expression tag	UNP P63086
A	2	MET	-	expression tag	UNP P63086

- Molecule 2 is a protein called DESIGNED ANKYRIN REPEAT PROTEIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	154	Total	C	N	O	S	0	0	0
			1167	728	206	232	1			

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	183	Total	O	0	0
			183	183		
3	B	53	Total	O	0	0
			53	53		

4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	68.04Å 89.37Å 99.54Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	19.43 – 1.97 19.43 – 1.97	Depositor EDS
% Data completeness (in resolution range)	93.9 (19.43-1.97) 93.7 (19.43-1.97)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.64 (at 1.97Å)	Xtriage
Refinement program	REFMAC 5.5.0109	Depositor
R, R_{free}	0.219 , 0.269 0.225 , 0.270	Depositor DCC
R_{free} test set	2038 reflections (4.98%)	wwPDB-VP
Wilson B-factor (Å ²)	27.5	Xtriage
Anisotropy	0.040	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 36.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.43$, $\langle L^2 \rangle = 0.26$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	4220	wwPDB-VP
Average B, all atoms (Å ²)	30.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.52% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

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	1.27	4/2884 (0.1%)	1.24	13/3907 (0.3%)
2	B	1.17	0/1187	1.14	1/1615 (0.1%)
All	All	1.25	4/4071 (0.1%)	1.21	14/5522 (0.3%)

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	A	0	1

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	144	LEU	C-O	6.15	1.31	1.24
1	A	20	GLY	N-CA	5.85	1.52	1.44
1	A	110	LEU	C-O	-5.21	1.17	1.24
1	A	215	ILE	N-CA	5.04	1.52	1.46

All (14) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	20	GLY	CA-C-N	-10.04	109.50	119.64
1	A	20	GLY	C-N-CA	-10.04	109.50	119.64
1	A	97	LYS	N-CA-C	-8.70	102.80	113.41
1	A	299	ARG	NE-CZ-NH2	-6.81	113.07	119.20
1	A	90	ALA	CA-C-N	6.59	126.28	119.82
1	A	90	ALA	C-N-CA	6.59	126.28	119.82
1	A	299	ARG	NE-CZ-NH1	5.74	127.24	121.50
1	A	188	THR	CB-CA-C	-5.55	99.68	109.62
1	A	185	TYR	N-CA-C	5.49	119.08	112.38
2	B	126	LEU	N-CA-C	5.49	117.97	111.33

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	57	PHE	N-CA-C	5.29	119.36	113.01
1	A	343	GLU	N-CA-C	5.12	116.55	111.07
1	A	98	ASP	N-CA-C	5.10	117.13	109.23
1	A	299	ARG	CD-NE-CZ	5.05	131.48	124.40

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	35	GLY	Peptide

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	2817	0	2820	94	0
2	B	1167	0	1128	19	0
3	A	183	0	0	9	0
3	B	53	0	0	2	0
All	All	4220	0	3948	106	0

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 13.

All (106) 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)
2:B:163:GLU:HG2	3:B:2051:HOH:O	1.56	1.04
1:A:229:LYS:NZ	2:B:79:ASP:OD2	2.03	0.91
2:B:153:SER:HB3	2:B:161:LEU:HD23	1.52	0.91
1:A:92:THR:HG23	1:A:95:GLN:HE21	1.40	0.86
1:A:78:HIS:HD2	1:A:80:ASN:H	1.28	0.81
1:A:174:PRO:O	1:A:178:HIS:CE1	2.38	0.77
1:A:89:ARG:HB2	1:A:96:MET:HE2	1.68	0.75
1:A:78:HIS:CD2	1:A:80:ASN:H	2.07	0.72

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:46:GLN:O	2:B:46:GLN:HG3	1.87	0.72
1:A:230:HIS:HE1	2:B:110:ASP:OD2	1.72	0.72
1:A:270:LYS:HE2	3:A:2151:HOH:O	1.88	0.72
1:A:92:THR:OG1	1:A:95:GLN:HG3	1.91	0.69
1:A:154:LEU:HD11	1:A:164:CYS:SG	2.35	0.67
1:A:168:LEU:HD21	1:A:184:GLU:HG2	1.78	0.65
1:A:153:LEU:CD2	1:A:163:ILE:HG12	2.27	0.64
1:A:82:ILE:HG21	1:A:164:CYS:HB2	1.79	0.64
1:A:230:HIS:CD2	1:A:233:ASP:H	2.18	0.62
1:A:176:HIS:O	1:A:178:HIS:N	2.32	0.61
1:A:118:HIS:HE1	3:A:2071:HOH:O	1.84	0.60
1:A:92:THR:CG2	1:A:95:GLN:HE21	2.13	0.59
1:A:295:ASN:HD22	1:A:295:ASN:C	2.11	0.59
1:A:90:ALA:HB1	1:A:95:GLN:HB2	1.87	0.57
1:A:164:CYS:O	1:A:165:ASP:CB	2.52	0.56
1:A:170:ARG:CZ	1:A:179:THR:HG22	2.35	0.56
1:A:197:MET:SD	1:A:234:GLN:HG2	2.45	0.56
1:A:89:ARG:HD3	1:A:96:MET:HE2	1.87	0.56
1:A:146:ARG:HG2	1:A:207:ILE:HD11	1.87	0.56
1:A:10:GLU:HA	1:A:10:GLU:OE1	2.07	0.55
1:A:168:LEU:HD11	1:A:184:GLU:HG3	1.88	0.55
1:A:289:ASP:OD2	3:A:2161:HOH:O	2.17	0.55
1:A:293:THR:O	1:A:299:ARG:HD2	2.07	0.55
1:A:31:GLU:HG2	1:A:32:GLY:N	2.23	0.54
1:A:239:LEU:HD22	1:A:245:PRO:HD3	1.89	0.54
1:A:178:HIS:O	1:A:179:THR:C	2.51	0.54
1:A:35:GLY:HA2	1:A:53:LYS:O	2.09	0.52
1:A:295:ASN:HD22	1:A:296:PRO:HD2	1.75	0.52
1:A:197:MET:CE	1:A:234:GLN:HG2	2.40	0.51
1:A:28:TYR:OH	1:A:31:GLU:HB3	2.11	0.51
2:B:114:LEU:HD22	2:B:143:ASP:HB2	1.92	0.51
1:A:89:ARG:CD	1:A:96:MET:HE2	2.41	0.50
1:A:295:ASN:ND2	1:A:297:HIS:H	2.09	0.50
1:A:55:SER:HB3	1:A:98:ASP:OD1	2.12	0.50
1:A:328:LYS:HE3	1:A:328:LYS:HA	1.92	0.49
1:A:230:HIS:CE1	2:B:110:ASP:OD2	2.60	0.49
1:A:177:ASP:O	1:A:177:ASP:OD1	2.30	0.49
1:A:89:ARG:CB	1:A:96:MET:HE2	2.37	0.49
1:A:295:ASN:HD22	1:A:296:PRO:N	2.10	0.49
1:A:301:GLU:H	1:A:304:GLN:HE21	1.60	0.49
1:A:178:HIS:O	1:A:179:THR:O	2.30	0.49

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:188:THR:HG22	1:A:190:TRP:H	1.77	0.49
1:A:195:GLU:O	1:A:200:SER:HB2	2.13	0.49
1:A:174:PRO:HD3	3:A:2087:HOH:O	2.13	0.49
1:A:295:ASN:ND2	1:A:296:PRO:HD2	2.27	0.48
1:A:205:LYS:HG3	3:A:2120:HOH:O	2.12	0.48
1:A:295:ASN:HD22	1:A:296:PRO:CD	2.26	0.48
1:A:89:ARG:HD3	1:A:96:MET:CE	2.44	0.48
1:A:174:PRO:O	1:A:178:HIS:NE2	2.45	0.48
1:A:229:LYS:HZ1	2:B:46:GLN:HE22	1.62	0.48
2:B:58:GLY:O	2:B:60:PRO:HD3	2.14	0.48
1:A:189:ARG:HD3	1:A:192:ARG:NH1	2.29	0.48
1:A:154:LEU:CD1	1:A:164:CYS:SG	3.01	0.48
1:A:78:HIS:HE1	3:A:2077:HOH:O	1.96	0.47
2:B:166:GLN:C	3:B:2053:HOH:O	2.57	0.47
1:A:178:HIS:HB2	3:A:2102:HOH:O	2.15	0.47
1:A:301:GLU:H	1:A:304:GLN:NE2	2.12	0.47
1:A:111:TYR:CE1	1:A:115:LYS:HE3	2.50	0.47
1:A:89:ARG:CG	1:A:96:MET:HE2	2.43	0.47
2:B:26:GLN:O	2:B:30:VAL:HG23	2.14	0.47
1:A:176:HIS:O	1:A:177:ASP:C	2.58	0.47
1:A:293:THR:O	1:A:299:ARG:CD	2.63	0.47
1:A:112:LYS:O	1:A:116:THR:HG23	2.14	0.47
1:A:168:LEU:HD11	1:A:184:GLU:CG	2.44	0.46
1:A:82:ILE:O	1:A:82:ILE:HG23	2.14	0.46
1:A:88:ILE:HB	1:A:100:TYR:HB2	1.97	0.46
1:A:82:ILE:HB	1:A:164:CYS:HB3	1.98	0.45
1:A:229:LYS:NZ	2:B:46:GLN:HE22	2.14	0.45
2:B:130:GLU:OE2	2:B:164:ILE:HG21	2.17	0.44
1:A:201:LYS:HD3	1:A:201:LYS:HA	1.30	0.44
2:B:130:GLU:OE2	2:B:164:ILE:CG2	2.65	0.44
1:A:113:LEU:HD13	1:A:119:LEU:HD21	2.00	0.44
1:A:198:LEU:HD13	1:A:235:LEU:HD21	1.99	0.44
1:A:229:LYS:HZ2	2:B:79:ASP:CG	2.15	0.44
1:A:153:LEU:HD22	1:A:163:ILE:HG12	1.99	0.43
1:A:67:LEU:HG	1:A:345:ILE:HG13	2.01	0.43
1:A:197:MET:HG3	1:A:231:TYR:HB2	1.98	0.43
1:A:198:LEU:HD21	1:A:262:LEU:HD11	2.00	0.43
2:B:37:GLY:O	2:B:38:ALA:C	2.59	0.43
1:A:111:TYR:HE1	1:A:115:LYS:HE3	1.84	0.42
1:A:252:CYS:HB3	3:A:2146:HOH:O	2.18	0.42
1:A:82:ILE:HG21	1:A:164:CYS:CB	2.46	0.42

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:197:MET:HE2	1:A:197:MET:HB2	1.78	0.42
1:A:232:LEU:HD13	2:B:144:LYS:HG2	2.01	0.42
1:A:56:PRO:HD2	1:A:57:PHE:CE2	2.54	0.42
1:A:89:ARG:CD	1:A:96:MET:CE	2.98	0.41
2:B:34:MET:CE	2:B:69:HIS:CG	3.04	0.41
1:A:57:PHE:CD1	1:A:342:LYS:HG3	2.55	0.41
1:A:94:GLU:H	1:A:94:GLU:CD	2.27	0.41
1:A:59:HIS:HB3	1:A:62:TYR:CD2	2.56	0.41
1:A:63:CYS:O	1:A:64:GLN:C	2.63	0.41
1:A:82:ILE:CG2	1:A:164:CYS:CB	2.99	0.41
1:A:267:HIS:HE1	1:A:269:ASN:OD1	2.04	0.41
2:B:24:ALA:CB	2:B:26:GLN:HE21	2.34	0.41
1:A:324:GLU:HG2	3:A:2080:HOH:O	2.20	0.40
1:A:170:ARG:NH1	1:A:330:ASP:OD1	2.53	0.40
1:A:257:LYS:HB3	1:A:257:LYS:HE2	1.79	0.40
1:A:144:LEU:HD22	1:A:205:LYS:HA	2.04	0.40

There are no symmetry-related clashes.

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	342/365 (94%)	318 (93%)	19 (6%)	5 (2%)	8	2
2	B	152/169 (90%)	147 (97%)	5 (3%)	0	100	100
All	All	494/534 (92%)	465 (94%)	24 (5%)	5 (1%)	12	5

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	177	ASP
1	A	179	THR

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	A	165	ASP
1	A	56	PRO
1	A	201	LYS

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	312/324 (96%)	282 (90%)	30 (10%)	8	2
2	B	119/132 (90%)	113 (95%)	6 (5%)	22	10
All	All	431/456 (94%)	395 (92%)	36 (8%)	10	3

All (36) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	10	GLU
1	A	34	TYR
1	A	59	HIS
1	A	67	LEU
1	A	73	LEU
1	A	74	LEU
1	A	88	ILE
1	A	92	THR
1	A	93	ILE
1	A	113	LEU
1	A	148	LEU
1	A	176	HIS
1	A	178	HIS
1	A	181	PHE
1	A	188	THR
1	A	197	MET
1	A	198	LEU
1	A	199	ASN
1	A	200	SER
1	A	201	LYS

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	A	235	LEU
1	A	250	LEU
1	A	253	ILE
1	A	256	LEU
1	A	293	THR
1	A	295	ASN
1	A	299	ARG
1	A	313	GLN
1	A	334	ASP
1	A	340	LYS
2	B	46	GLN
2	B	68	LYS
2	B	114	LEU
2	B	126	LEU
2	B	130	GLU
2	B	166	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (22) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	15	GLN
1	A	60	GLN
1	A	78	HIS
1	A	85	ASN
1	A	95	GLN
1	A	118	HIS
1	A	176	HIS
1	A	199	ASN
1	A	230	HIS
1	A	234	GLN
1	A	251	ASN
1	A	267	HIS
1	A	295	ASN
1	A	297	HIS
1	A	304	GLN
2	B	26	GLN
2	B	36	ASN
2	B	46	GLN
2	B	59	HIS
2	B	69	HIS
2	B	156	ASN
2	B	166	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 oligosaccharides 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

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, 95th 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 ²)	Q<0.9
1	A	344/365 (94%)	0.32	38 (11%) 10 14	13, 25, 59, 81	0
2	B	154/169 (91%)	0.23	7 (4%) 38 47	17, 29, 52, 68	0
All	All	498/534 (93%)	0.30	45 (9%) 15 21	13, 27, 57, 81	0

All (45) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	34	TYR	5.1
1	A	200	SER	4.2
1	A	177	ASP	4.1
1	A	164	CYS	4.0
1	A	176	HIS	3.9
1	A	337	PRO	3.8
1	A	32	GLY	3.5
2	B	165	LEU	3.4
1	A	175	ASP	3.3
1	A	90	ALA	3.3
1	A	182	LEU	3.3
1	A	179	THR	3.2
2	B	24	ALA	3.2
1	A	336	LEU	3.1
1	A	181	PHE	3.1
1	A	163	ILE	3.0
1	A	174	PRO	2.9
1	A	334	ASP	2.9
1	A	62	TYR	2.8
2	B	164	ILE	2.8
1	A	178	HIS	2.7
1	A	59	HIS	2.7
1	A	61	THR	2.6
1	A	99	VAL	2.6

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	A	183	THR	2.6
1	A	186	VAL	2.6
1	A	64	GLN	2.4
1	A	33	ALA	2.4
1	A	201	LYS	2.4
2	B	56	TRP	2.4
1	A	346	PHE	2.4
2	B	166	GLN	2.4
1	A	333	LEU	2.3
1	A	184	GLU	2.3
1	A	203	TYR	2.3
1	A	91	PRO	2.3
2	B	163	GLU	2.2
1	A	180	GLY	2.2
1	A	341	LEU	2.2
2	B	157	GLY	2.2
1	A	172	ALA	2.2
1	A	97	LYS	2.2
1	A	20	GLY	2.1
1	A	57	PHE	2.1
1	A	56	PRO	2.0

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 oligosaccharides in this entry.

6.4 Ligands [i](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.