



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 05:31 PM UTC

PDB ID : 6VTC / pdb_00006vtc
Title : p53-specific T cell receptor
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Deposited on : 2020-02-12
Resolution : 1.83 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

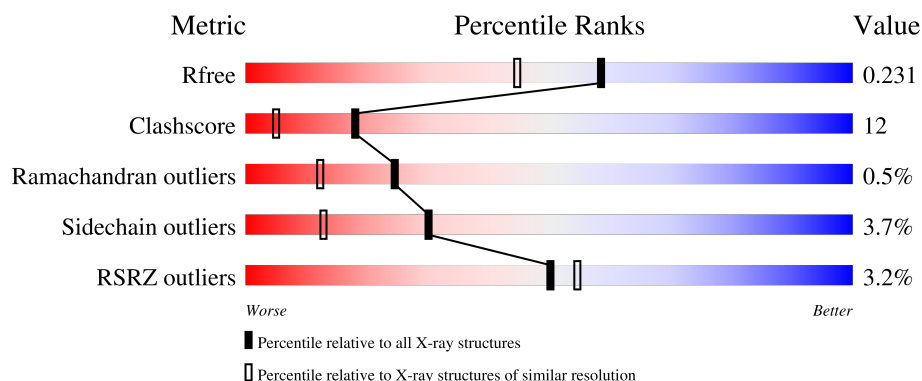
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

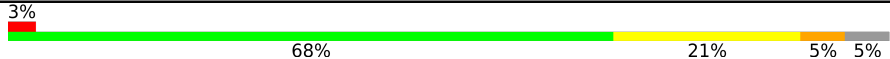
The reported resolution of this entry is 1.83 Å.

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	1296 (1.84-1.84)
Clashscore	190562	1329 (1.84-1.84)
Ramachandran outliers	187476	1318 (1.84-1.84)
Sidechain outliers	187428	1318 (1.84-1.84)
RSRZ outliers	180081	1296 (1.84-1.84)

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	208	
2	B	244	

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 3824 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 T-cell Receptor 1a2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	197	Total	C	N	O	S	0	0	0
			1513	950	242	311	10			

- Molecule 2 is a protein called p53-specific T cell receptor, B-chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	243	Total	C	N	O	S	0	0	0
			1920	1213	334	365	8			

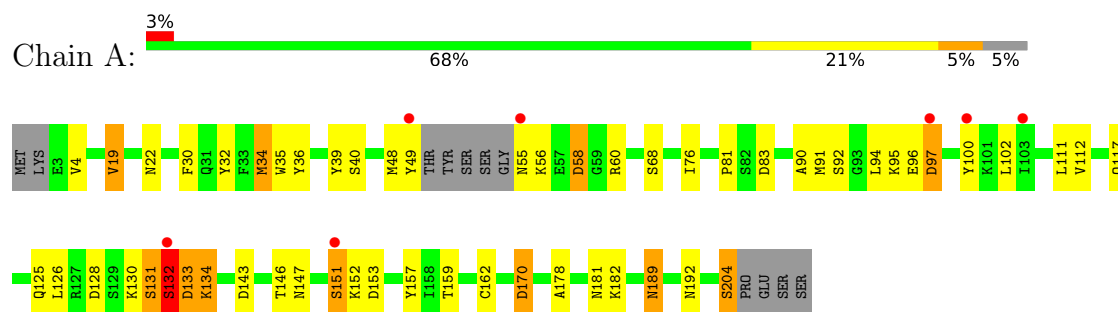
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	144	Total	O	0	0
			144	144		
3	B	247	Total	O	0	0
			247	247		

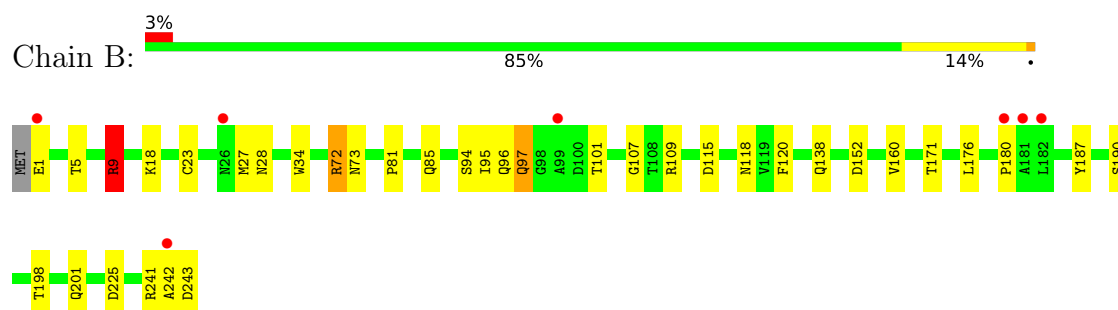
3 Residue-property plots [i](#)

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.

• Molecule 1: T-cell Receptor 1a2



• Molecule 2: p53-specific T cell receptor, B-chain



4 Data and refinement statistics

Property	Value	Source
Space group	P 2 ₁ 2 ₁ 2 ₁	Depositor
Cell constants a, b, c, α , β , γ	41.12Å 106.37Å 112.69Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.00 – 1.83 20.00 – 1.83	Depositor EDS
% Data completeness (in resolution range)	99.7 (20.00-1.83) 99.7 (20.00-1.83)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.73 (at 1.83Å)	Xtriage
Refinement program	REFMAC 5.8.0238	Depositor
R, R_{free}	0.190 , 0.231 0.190 , 0.231	Depositor DCC
R_{free} test set	2174 reflections (4.88%)	wwPDB-VP
Wilson B-factor (Å ²)	34.6	Xtriage
Anisotropy	0.452	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 47.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	3824	wwPDB-VP
Average B, all atoms (Å ²)	43.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.56% 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	0.90	0/1546	1.44	13/2098 (0.6%)
2	B	0.99	1/1972 (0.1%)	1.43	14/2691 (0.5%)
All	All	0.95	1/3518 (0.0%)	1.43	27/4789 (0.6%)

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
2	B	0	1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	72	ARG	NE-CZ	6.11	1.39	1.33

All (27) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	5	THR	CA-CB-OG1	-8.38	97.03	109.60
1	A	162	CYS	N-CA-CB	-7.41	97.72	111.13
2	B	120	PHE	CB-CA-C	-7.13	99.11	109.45
1	A	97	ASP	CA-CB-CG	6.49	119.09	112.60
2	B	72	ARG	NE-CZ-NH2	6.49	125.04	119.20
1	A	128	ASP	CA-CB-CG	6.28	118.88	112.60
2	B	198	THR	CA-CB-OG1	-6.24	100.24	109.60
1	A	147	ASN	CA-C-N	-6.21	114.95	122.90
1	A	147	ASN	C-N-CA	-6.21	114.95	122.90
2	B	101	THR	CA-CB-OG1	-6.01	100.59	109.60
2	B	160	VAL	CA-C-O	-5.87	114.29	120.39
2	B	180	PRO	N-CA-C	-5.82	106.88	114.03
1	A	133	ASP	CA-CB-CG	5.73	118.33	112.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	72	ARG	CD-NE-CZ	5.59	132.23	124.40
2	B	187	TYR	CB-CA-C	-5.57	98.37	109.79
1	A	143	ASP	CA-CB-CG	5.38	117.97	112.60
1	A	134	LYS	CA-C-N	-5.36	113.79	122.36
1	A	134	LYS	C-N-CA	-5.36	113.79	122.36
2	B	115	ASP	CA-CB-CG	-5.27	107.33	112.60
2	B	9	ARG	CB-CG-CD	5.23	123.34	111.30
1	A	58	ASP	CA-CB-CG	5.22	117.82	112.60
2	B	171	THR	CA-CB-OG1	-5.13	101.90	109.60
1	A	170	ASP	CB-CA-C	5.12	119.97	112.09
2	B	118	ASN	CA-CB-CG	-5.09	107.50	112.60
2	B	97	GLN	CB-CG-CD	-5.08	103.95	112.60
1	A	132	SER	CA-C-N	-5.03	115.88	124.82
1	A	132	SER	C-N-CA	-5.03	115.88	124.82

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	B	1	GLU	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	1513	0	1392	65	0
2	B	1920	0	1824	18	0
3	A	144	0	0	8	0
3	B	247	0	0	9	0
All	All	3824	0	3216	80	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 12.

All (80) 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:A:4:VAL:HG22	1:A:91:MET:CE	1.57	1.34
1:A:4:VAL:CG2	1:A:91:MET:HE2	1.77	1.12
1:A:4:VAL:HG22	1:A:91:MET:HE2	1.21	1.09
1:A:4:VAL:HG22	1:A:91:MET:HE1	1.36	1.07
1:A:34:MET:HE1	3:B:495:HOH:O	1.53	1.05
1:A:130:LYS:O	1:A:131:SER:HB2	1.62	0.96
1:A:34:MET:CE	3:B:495:HOH:O	2.13	0.94
1:A:95:LYS:HB2	3:A:305:HOH:O	1.68	0.93
1:A:30:PHE:CE2	1:A:91:MET:HE3	2.07	0.90
1:A:32:TYR:CE1	1:A:94:LEU:HD23	2.14	0.81
1:A:49:TYR:O	1:A:55:ASN:N	2.16	0.79
1:A:34:MET:HG2	1:A:90:ALA:HB3	1.63	0.79
1:A:151:SER:OG	3:A:301:HOH:O	2.00	0.79
1:A:32:TYR:HD2	3:B:534:HOH:O	1.64	0.79
1:A:48:MET:HE1	1:A:58:ASP:HB3	1.66	0.77
1:A:32:TYR:HE1	1:A:94:LEU:HD23	1.51	0.75
1:A:32:TYR:CE1	1:A:94:LEU:CD2	2.70	0.74
1:A:32:TYR:HE1	1:A:94:LEU:CD2	1.99	0.74
1:A:204:SER:O	3:A:303:HOH:O	2.05	0.74
1:A:132:SER:O	1:A:133:ASP:CB	2.35	0.74
1:A:100:TYR:H	2:B:97:GLN:HE21	1.38	0.71
1:A:56:LYS:HG2	3:A:432:HOH:O	1.91	0.70
2:B:9:ARG:NH2	2:B:107:GLY:O	2.25	0.69
2:B:109:ARG:NH2	2:B:152:ASP:HB3	2.08	0.69
1:A:4:VAL:CG2	1:A:91:MET:CE	2.44	0.68
1:A:151:SER:CB	3:A:301:HOH:O	2.42	0.67
1:A:30:PHE:HE2	1:A:91:MET:HE3	1.55	0.65
1:A:132:SER:O	1:A:133:ASP:HB2	1.96	0.64
1:A:130:LYS:O	1:A:131:SER:CB	2.39	0.61
1:A:60:ARG:NH2	1:A:83:ASP:OD2	2.32	0.60
2:B:94:SER:OG	2:B:96:GLN:O	2.20	0.60
1:A:132:SER:HB2	1:A:134:LYS:HG3	1.84	0.59
2:B:18:LYS:NZ	3:B:303:HOH:O	2.34	0.59
1:A:181:ASN:N	3:A:302:HOH:O	2.04	0.59
1:A:19:VAL:HG13	1:A:76:ILE:HB	1.86	0.57
2:B:109:ARG:HH22	2:B:152:ASP:HB3	1.68	0.57
1:A:34:MET:HB3	1:A:49:TYR:HA	1.88	0.56
1:A:94:LEU:O	1:A:97:ASP:HB2	2.06	0.56
1:A:34:MET:CG	1:A:90:ALA:HB3	2.35	0.55
1:A:35:TRP:N	1:A:48:MET:O	2.36	0.54
2:B:72:ARG:NH2	2:B:73:ASN:HD21	2.07	0.53
1:A:34:MET:HE1	3:B:534:HOH:O	2.09	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:131:SER:O	1:A:132:SER:CB	2.56	0.52
1:A:32:TYR:CE1	1:A:94:LEU:HD22	2.45	0.52
1:A:152:LYS:HA	3:A:375:HOH:O	2.10	0.52
1:A:4:VAL:HG21	1:A:91:MET:HE2	1.83	0.51
1:A:55:ASN:CG	1:A:56:LYS:H	2.18	0.50
1:A:100:TYR:H	2:B:97:GLN:NE2	2.06	0.50
2:B:225:ASP:N	3:B:310:HOH:O	2.44	0.50
1:A:131:SER:O	1:A:132:SER:HB3	2.12	0.49
1:A:92:SER:HB3	1:A:102:LEU:HA	1.96	0.48
1:A:96:GLU:OE1	3:A:304:HOH:O	2.20	0.48
1:A:132:SER:CB	1:A:134:LYS:HG3	2.44	0.48
2:B:242:ALA:O	2:B:243:ASP:C	2.58	0.47
2:B:138:GLN:NE2	3:B:302:HOH:O	2.32	0.46
1:A:159:THR:HG21	2:B:190:SER:OG	2.16	0.46
2:B:201:GLN:HA	2:B:241:ARG:O	2.16	0.45
1:A:32:TYR:HE1	1:A:94:LEU:HA	1.80	0.45
1:A:126:LEU:HD12	1:A:126:LEU:N	2.31	0.45
1:A:32:TYR:CE1	1:A:94:LEU:HA	2.52	0.45
1:A:34:MET:HB3	1:A:34:MET:HE3	1.62	0.45
1:A:34:MET:HE2	3:B:485:HOH:O	2.17	0.45
2:B:81:PRO:HA	2:B:85:GLN:OE1	2.17	0.45
1:A:125:GLN:C	1:A:126:LEU:HD12	2.41	0.45
2:B:95:ILE:HG22	2:B:96:GLN:HG3	1.99	0.45
1:A:48:MET:HE1	1:A:58:ASP:CB	2.43	0.44
1:A:181:ASN:O	1:A:182:LYS:C	2.60	0.43
1:A:153:ASP:OD1	1:A:153:ASP:C	2.61	0.43
1:A:39:TYR:O	1:A:40:SER:C	2.62	0.42
1:A:34:MET:CB	1:A:49:TYR:HA	2.49	0.42
1:A:19:VAL:CG1	1:A:76:ILE:HB	2.48	0.42
2:B:138:GLN:HG2	3:B:493:HOH:O	2.19	0.42
1:A:132:SER:O	1:A:133:ASP:HB3	2.19	0.42
1:A:32:TYR:HE1	1:A:94:LEU:HD22	1.77	0.42
1:A:189:ASN:HA	1:A:192:ASN:OD1	2.20	0.42
2:B:23:CYS:HB2	2:B:34:TRP:CZ2	2.55	0.42
1:A:157:TYR:O	1:A:178:ALA:HA	2.20	0.41
2:B:176:LEU:C	2:B:176:LEU:HD12	2.45	0.41
1:A:34:MET:HG3	1:A:36:TYR:HE1	1.85	0.41
1:A:81:PRO:HA	1:A:112:VAL:HB	2.02	0.41

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	193/208 (93%)	184 (95%)	7 (4%)	2 (1%)	12	4
2	B	241/244 (99%)	236 (98%)	5 (2%)	0	100	100
All	All	434/452 (96%)	420 (97%)	12 (3%)	2 (0%)	24	12

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	132	SER
1	A	131	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	A	169/189 (89%)	158 (94%)	11 (6%)	15	3
2	B	208/217 (96%)	205 (99%)	3 (1%)	59	45
All	All	377/406 (93%)	363 (96%)	14 (4%)	30	13

All (14) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	19	VAL
1	A	22	ASN
1	A	34	MET
1	A	68	SER

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Mol	Chain	Res	Type
1	A	111	LEU
1	A	117	GLN
1	A	146	THR
1	A	151	SER
1	A	170	ASP
1	A	189	ASN
1	A	204	SER
2	B	9	ARG
2	B	27	MET
2	B	28	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	80	GLN
1	A	189	ASN
2	B	73	ASN
2	B	96	GLN
2	B	97	GLN
2	B	179	GLN
2	B	183	ASN

5.3.3 RNA ⓘ

There are no RNA molecules 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 oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

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](#)

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	197/208 (94%)	0.21	7 (3%) 46 49	27, 43, 77, 97	0
2	B	243/244 (99%)	-0.15	7 (2%) 53 58	25, 35, 59, 91	0
All	All	440/452 (97%)	0.01	14 (3%) 50 54	25, 39, 70, 97	0

All (14) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	100	TYR	3.8
2	B	181	ALA	3.6
1	A	97	ASP	2.5
2	B	182	LEU	2.5
1	A	55	ASN	2.4
1	A	49	TYR	2.4
2	B	1	GLU	2.4
1	A	103	ILE	2.3
2	B	180	PRO	2.3
2	B	99	ALA	2.2
2	B	26	ASN	2.2
2	B	242	ALA	2.1
1	A	132	SER	2.0
1	A	151	SER	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.