



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 12, 2026 – 09:54 AM UTC

PDB ID : 1XJV / pdb_00001xjv
Title : Crystal structure of human POT1 bound to telomeric single-stranded DNA (TTAGGGTTAG)
Authors : Lei, M.; Podell, E.R.; Cech, T.R.
Deposited on : 2004-09-25
Resolution : 1.73 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

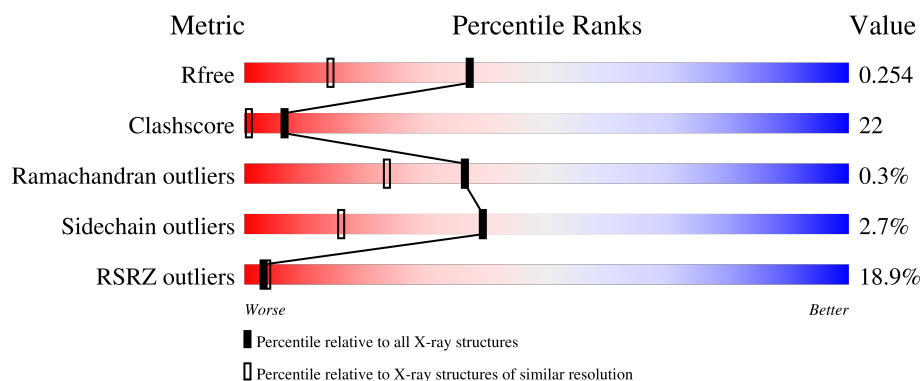
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.73 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	1187 (1.74-1.74)
Clashscore	190562	1207 (1.74-1.74)
Ramachandran outliers	187476	1200 (1.74-1.74)
Sidechain outliers	187428	1200 (1.74-1.74)
RSRZ outliers	180081	1188 (1.74-1.74)

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	B	10	50% 20% 70% 10%
2	A	294	18% 73% 22% . .

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 2766 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 DNA chain called hT10 d(TTAGGGTTAG).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	B	10	Total	C	N	O	P	0	0	0
			207	100	38	60	9			

- Molecule 2 is a protein called Protection of telomeres 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	A	291	Total	C	N	O	S	0	0	0
			2312	1484	390	429	9			

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	15	Total	O	0	0
			15	15		
3	A	232	Total	O	0	0
			232	232		

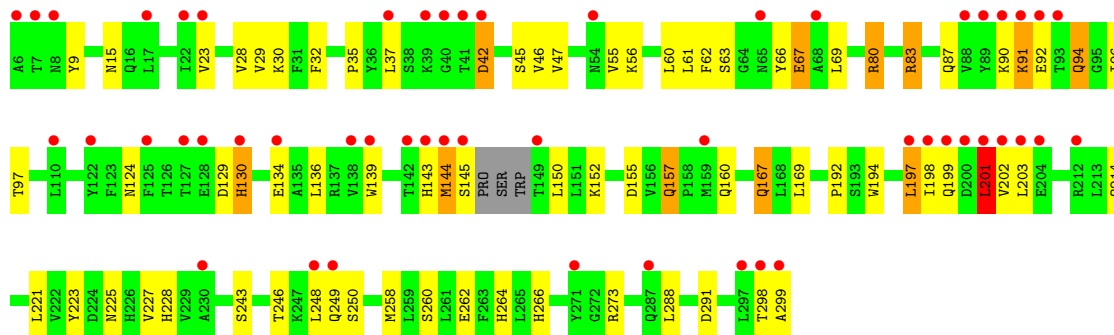
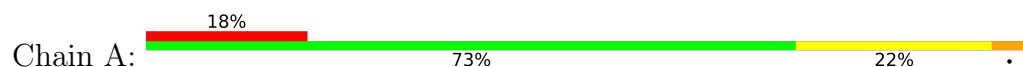
3 Residue-property plots [i](#)

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: hT10 d(TTAGGGTTAG)



- Molecule 2: Protection of telomeres 1



4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	102.03Å 103.24Å 71.66Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	36.29 – 1.73 36.29 – 1.73	Depositor EDS
% Data completeness (in resolution range)	(Not available) (36.29-1.73) 92.4 (36.29-1.73)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.04	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.78 (at 1.73Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.228 , 0.237 0.239 , 0.254	Depositor DCC
R_{free} test set	180 reflections (0.47%)	wwPDB-VP
Wilson B-factor (Å ²)	26.6	Xtriage
Anisotropy	0.431	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 37.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.025 for -k,-h,-l 0.014 for -1/2*h-1/2*k+l,-1/2*h-1/2*k-l,1/2 *h-1/2*k 0.013 for -1/2*h-1/2*k-l,-1/2*h-1/2*k+l,-1/ 2*h+1/2*k 0.015 for -1/2*h+1/2*k-l,1/2*h-1/2*k-l,-1/2 *h-1/2*k 0.000 for -1/2*h+1/2*k+l,1/2*h-1/2*k+l,1/ 2*h+1/2*k	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	2766	wwPDB-VP
Average B, all atoms (Å ²)	37.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.83% of the height of the origin peak. No significant pseudotranslation is detected.*

¹ Intensities estimated from amplitudes.

² Theoretical 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 [i](#)

5.1 Standard geometry [i](#)

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	0.28	0/232	0.83	1/358 (0.3%)
2	A	0.48	0/2362	1.01	13/3202 (0.4%)
All	All	0.46	0/2594	0.99	14/3560 (0.4%)

There are no bond length outliers.

All (14) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	227	VAL	N-CA-C	7.07	117.20	110.42
2	A	201	LEU	CA-C-N	7.03	129.56	120.56
2	A	201	LEU	C-N-CA	7.03	129.56	120.56
2	A	214	GLN	CB-CA-C	-6.25	109.36	116.54
1	B	6	DG	C2'-C3'-O3'	6.10	120.65	111.50
2	A	194	TRP	N-CA-C	-5.83	103.24	110.41
2	A	23	VAL	N-CA-C	5.75	115.83	108.82
2	A	197	LEU	N-CA-CB	5.58	118.31	109.71
2	A	167	GLN	OE1-CD-NE2	-5.42	117.18	122.60
2	A	202	VAL	N-CA-C	5.37	115.58	110.53
2	A	157	GLN	CA-C-N	5.16	126.29	119.84
2	A	157	GLN	C-N-CA	5.16	126.29	119.84
2	A	42	ASP	N-CA-C	5.07	117.97	110.46
2	A	201	LEU	N-CA-CB	-5.05	103.28	110.80

There are no chirality outliers.

There are no planarity outliers.

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	B	207	0	116	22	0
2	A	2312	0	2329	89	3
3	A	232	0	0	35	3
3	B	15	0	0	2	0
All	All	2766	0	2445	107	3

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 22.

All (107) 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:A:28:VAL:HG11	2:A:144:MET:HE1	1.08	1.05
2:A:28:VAL:CG1	2:A:144:MET:HE1	1.85	1.04
2:A:144:MET:HG2	2:A:144:MET:O	1.59	1.02
2:A:198:ILE:HD11	2:A:203:LEU:HB2	1.41	1.02
2:A:201:LEU:HD23	2:A:201:LEU:O	1.62	0.99
2:A:69:LEU:HD11	3:A:520:HOH:O	1.67	0.94
1:B:4:DG:H21	2:A:94:GLN:HE22	1.07	0.94
2:A:144:MET:O	2:A:145:SER:HB2	1.66	0.94
2:A:28:VAL:HG11	2:A:144:MET:CE	2.01	0.90
2:A:198:ILE:CD1	2:A:203:LEU:HB2	2.02	0.90
2:A:144:MET:O	2:A:145:SER:CB	2.18	0.90
2:A:87:GLN:HE21	2:A:94:GLN:HE21	1.21	0.89
1:B:4:DG:H3'	1:B:5:DG:H5''	1.57	0.86
2:A:29:VAL:HG13	3:A:525:HOH:O	1.75	0.86
2:A:45:SER:HB3	3:A:461:HOH:O	1.74	0.85
2:A:157:GLN:H	2:A:160:GLN:HE21	1.21	0.85
1:B:4:DG:C3'	1:B:5:DG:H5''	2.11	0.81
2:A:288:LEU:HD12	3:A:491:HOH:O	1.81	0.80
2:A:61:LEU:HB2	3:A:461:HOH:O	1.81	0.80
2:A:143:HIS:O	2:A:144:MET:HB2	1.80	0.78
2:A:291:ASP:HB2	3:A:491:HOH:O	1.84	0.78
2:A:69:LEU:HD22	3:A:461:HOH:O	1.84	0.76
2:A:169:LEU:HG	3:A:496:HOH:O	1.86	0.76
2:A:62:PHE:HA	3:A:520:HOH:O	1.86	0.75
2:A:87:GLN:HE21	2:A:94:GLN:NE2	1.84	0.73
1:B:8:DT:C2	2:A:266:HIS:CE1	2.78	0.71
2:A:288:LEU:HA	3:A:491:HOH:O	1.90	0.71
1:B:4:DG:H21	2:A:94:GLN:NE2	1.87	0.70
2:A:246:THR:HG23	3:A:459:HOH:O	1.91	0.70
2:A:298:THR:HG22	2:A:298:THR:O	1.92	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:6:DG:H2''	1:B:7:DT:OP2	1.92	0.70
1:B:8:DT:H2''	1:B:9:DA:O5'	1.92	0.69
2:A:197:LEU:HD12	2:A:198:ILE:H	1.57	0.69
2:A:143:HIS:O	2:A:144:MET:CB	2.41	0.69
2:A:9:TYR:HB3	3:A:526:HOH:O	1.93	0.68
2:A:87:GLN:NE2	2:A:94:GLN:HE21	1.92	0.67
2:A:260:SER:HA	3:A:468:HOH:O	1.95	0.67
2:A:152:LYS:HB2	3:A:462:HOH:O	1.95	0.66
2:A:243:SER:HB3	2:A:273:ARG:HG2	1.79	0.65
2:A:80:ARG:HG2	3:A:510:HOH:O	1.96	0.64
2:A:67:GLU:H	2:A:67:GLU:CD	2.06	0.64
2:A:197:LEU:HD12	2:A:198:ILE:N	2.14	0.62
2:A:201:LEU:HD23	2:A:201:LEU:C	2.24	0.62
2:A:225:ASN:HB2	3:A:492:HOH:O	1.99	0.62
2:A:150:LEU:HA	3:A:515:HOH:O	2.01	0.60
2:A:167:GLN:HG2	3:A:496:HOH:O	2.01	0.60
2:A:248:LEU:HD12	2:A:260:SER:O	2.02	0.60
2:A:42:ASP:HB3	2:A:63:SER:C	2.27	0.59
1:B:8:DT:H72	1:B:9:DA:C2	2.38	0.59
2:A:152:LYS:NZ	3:A:491:HOH:O	2.36	0.58
1:B:8:DT:O2	2:A:266:HIS:CE1	2.57	0.58
2:A:139:TRP:CE2	2:A:143:HIS:HB2	2.37	0.58
2:A:91:LYS:HB3	2:A:91:LYS:NZ	2.19	0.57
2:A:299:ALA:HB2	3:A:425:HOH:O	2.04	0.56
2:A:250:SER:HA	2:A:258:MET:O	2.05	0.56
2:A:223:TYR:OH	2:A:264:HIS:HD2	1.89	0.56
2:A:130:HIS:O	2:A:134:GLU:HG3	2.06	0.56
2:A:249:GLN:NE2	2:A:262:GLU:OE2	2.36	0.56
2:A:155:ASP:N	3:A:462:HOH:O	2.38	0.55
2:A:35:PRO:HB2	2:A:66:TYR:CE1	2.42	0.55
2:A:42:ASP:HB3	2:A:63:SER:O	2.07	0.55
2:A:62:PHE:HD2	3:A:531:HOH:O	1.89	0.54
2:A:157:GLN:H	2:A:160:GLN:NE2	1.98	0.54
2:A:55:VAL:N	3:A:507:HOH:O	2.39	0.54
1:B:3:DA:H1'	3:B:173:HOH:O	2.07	0.53
1:B:4:DG:C3'	1:B:5:DG:C5'	2.86	0.53
1:B:6:DG:C2'	1:B:7:DT:OP2	2.57	0.52
1:B:8:DT:C2	1:B:9:DA:O4'	2.64	0.51
2:A:30:LYS:HE3	2:A:56:LYS:HE2	1.93	0.51
2:A:61:LEU:HA	2:A:97:THR:HG23	1.93	0.51
2:A:96:ILE:HG13	3:A:531:HOH:O	2.10	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:37:LEU:HB2	3:A:497:HOH:O	2.11	0.50
2:A:298:THR:O	2:A:299:ALA:HB2	2.11	0.50
2:A:60:LEU:HD11	2:A:94:GLN:OE1	2.12	0.49
2:A:9:TYR:OH	2:A:83:ARG:NH2	2.46	0.49
2:A:80:ARG:NH2	2:A:124:ASN:O	2.37	0.49
2:A:262:GLU:OE1	2:A:264:HIS:HE1	1.96	0.48
1:B:8:DT:O2	1:B:9:DA:H4'	2.13	0.48
1:B:4:DG:C5'	1:B:5:DG:H5''	2.45	0.47
2:A:228:HIS:HE1	3:A:492:HOH:O	1.98	0.47
1:B:8:DT:H1'	1:B:9:DA:H5'	1.97	0.47
2:A:90:LYS:C	2:A:92:GLU:H	2.23	0.47
1:B:8:DT:H2''	1:B:9:DA:O4'	2.15	0.46
2:A:97:THR:OG1	3:A:531:HOH:O	2.20	0.46
2:A:80:ARG:HH21	2:A:129:ASP:CG	2.23	0.46
2:A:198:ILE:HD13	2:A:203:LEU:HB2	1.95	0.46
2:A:62:PHE:HD1	3:A:457:HOH:O	1.99	0.46
1:B:2:DT:H3'	1:B:3:DA:C8	2.50	0.46
2:A:47:VAL:CG1	3:A:525:HOH:O	2.65	0.45
2:A:91:LYS:HB3	2:A:91:LYS:HZ2	1.80	0.45
2:A:221:LEU:HD23	2:A:223:TYR:OH	2.16	0.45
3:B:173:HOH:O	2:A:60:LEU:HD22	2.16	0.45
1:B:4:DG:C4'	1:B:5:DG:H5''	2.46	0.44
2:A:167:GLN:C	3:A:496:HOH:O	2.60	0.43
2:A:32:PHE:HA	2:A:46:VAL:O	2.19	0.43
1:B:2:DT:H5''	1:B:3:DA:N7	2.33	0.43
2:A:15:ASN:HB3	2:A:55:VAL:HG21	2.01	0.43
2:A:62:PHE:CA	3:A:520:HOH:O	2.57	0.43
2:A:223:TYR:OH	2:A:264:HIS:CD2	2.72	0.42
1:B:4:DG:H3'	1:B:5:DG:C5'	2.38	0.41
2:A:139:TRP:CD1	2:A:143:HIS:ND1	2.88	0.41
2:A:192:PRO:HB2	3:A:468:HOH:O	2.20	0.41
1:B:6:DG:N3	1:B:6:DG:H5''	2.36	0.41
2:A:30:LYS:HE3	2:A:56:LYS:CE	2.50	0.41
2:A:136:LEU:HD23	3:A:514:HOH:O	2.21	0.40
2:A:192:PRO:HD2	3:A:468:HOH:O	2.21	0.40
2:A:47:VAL:HG13	3:A:525:HOH:O	2.21	0.40

All (3) 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	Interatomic distance (Å)	Clash overlap (Å)
2:A:199:GLN:NE2	3:A:365:HOH:O[6_554]	1.68	0.52
2:A:199:GLN:CD	3:A:365:HOH:O[6_554]	1.78	0.42
2:A:199:GLN:OE1	3:A:365:HOH:O[6_554]	1.99	0.21

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
2	A	287/294 (98%)	276 (96%)	10 (4%)	1 (0%)	36 23

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	A	144	MET

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
2	A	261/264 (99%)	254 (97%)	7 (3%)	39 16

All (7) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	A	67	GLU
2	A	80	ARG
2	A	83	ARG

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Mol	Chain	Res	Type
2	A	91	LYS
2	A	94	GLN
2	A	130	HIS
2	A	201	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
2	A	8	ASN
2	A	24	ASN
2	A	65	ASN
2	A	75	ASN
2	A	94	GLN
2	A	160	GLN
2	A	245	HIS
2	A	255	ASN
2	A	264	HIS
2	A	266	HIS

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 ⓘ

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, 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	B	10/10 (100%)	2.49	5 (50%) 0 0	44, 59, 68, 71	0
2	A	291/294 (98%)	1.09	52 (17%) 3 4	20, 31, 55, 67	0
All	All	301/304 (99%)	1.14	57 (18%) 3 4	20, 32, 58, 71	0

All (57) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	A	7	THR	6.7
2	A	149	THR	6.3
2	A	200	ASP	5.8
2	A	198	ILE	5.2
2	A	299	ALA	5.2
1	B	8	DT	4.7
2	A	201	LEU	4.6
2	A	199	GLN	4.3
2	A	298	THR	4.2
2	A	6	ALA	4.1
2	A	142	THR	4.1
2	A	91	LYS	4.1
2	A	271	TYR	4.0
2	A	41	THR	4.0
2	A	145	SER	3.9
2	A	202	VAL	3.9
1	B	1	DT	3.8
2	A	125	PHE	3.7
1	B	9	DA	3.6
2	A	249	GLN	3.6
2	A	197	LEU	3.5
2	A	37	LEU	3.4
2	A	90	LYS	3.3
2	A	212	ARG	3.3

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Mol	Chain	Res	Type	RSRZ
2	A	143	HIS	3.2
2	A	144	MET	3.2
2	A	8	ASN	2.9
2	A	42	ASP	2.8
2	A	17	LEU	2.8
2	A	130	HIS	2.7
2	A	40	GLY	2.7
2	A	39	LYS	2.7
2	A	297	LEU	2.6
2	A	22	ILE	2.6
2	A	93	THR	2.5
1	B	2	DT	2.5
2	A	134	GLU	2.5
2	A	248	LEU	2.5
2	A	68	ALA	2.5
2	A	110	LEU	2.5
2	A	89	TYR	2.4
2	A	128	GLU	2.3
2	A	139	TRP	2.3
2	A	54	ASN	2.3
2	A	92	GLU	2.3
1	B	10	DG	2.3
2	A	127	THR	2.3
2	A	203	LEU	2.2
2	A	122	TYR	2.2
2	A	204	GLU	2.1
2	A	230	ALA	2.1
2	A	65	ASN	2.1
2	A	287	GLN	2.1
2	A	138	VAL	2.1
2	A	159	MET	2.0
2	A	23	VAL	2.0
2	A	88	VAL	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.