



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 12:37 PM GMT

PDB ID : 3RK6
Title : Crystal structure of the middle domain of human Paip1
Authors : Lei, J.; Mesters, J.R.; Hilgenfeld, R.
Deposited on : 2011-04-17
Resolution : 2.00 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
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467
Mogul : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

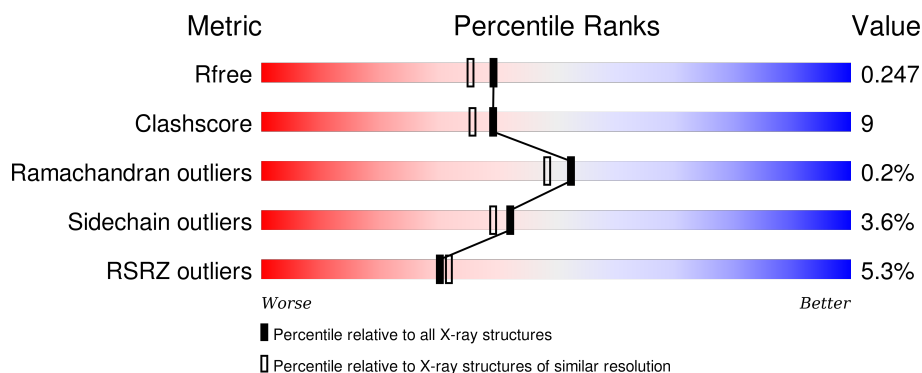
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION


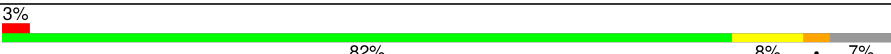
The reported resolution of this entry is 2.00 Å.

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))
R_{free}	91344	6249 (2.00-2.00)
Clashscore	102246	7340 (2.00-2.00)
Ramachandran outliers	100387	7248 (2.00-2.00)
Sidechain outliers	100360	7247 (2.00-2.00)
RSRZ outliers	91569	6262 (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 ≥ 3 , 2, 1 and 0 types of geometric quality criteria. 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	234	
1	B	234	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 3952 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 Polyadenylate-binding protein-interacting protein 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	217	Total	C	N	O	S	0	8	0
			1784	1126	301	346	11			
1	B	217	Total	C	N	O	S	0	10	0
			1798	1136	304	348	10			

There are 34 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	61	GLY	-	EXPRESSION TAG	UNP Q9H074
A	62	SER	-	EXPRESSION TAG	UNP Q9H074
A	63	HIS	-	EXPRESSION TAG	UNP Q9H074
A	64	MET	-	EXPRESSION TAG	UNP Q9H074
A	65	ALA	-	EXPRESSION TAG	UNP Q9H074
A	66	SER	-	EXPRESSION TAG	UNP Q9H074
A	67	MET	-	EXPRESSION TAG	UNP Q9H074
A	68	THR	-	EXPRESSION TAG	UNP Q9H074
A	69	GLY	-	EXPRESSION TAG	UNP Q9H074
A	70	GLY	-	EXPRESSION TAG	UNP Q9H074
A	71	GLN	-	EXPRESSION TAG	UNP Q9H074
A	72	GLN	-	EXPRESSION TAG	UNP Q9H074
A	73	MET	-	EXPRESSION TAG	UNP Q9H074
A	74	GLY	-	EXPRESSION TAG	UNP Q9H074
A	75	ARG	-	EXPRESSION TAG	UNP Q9H074
A	76	GLY	-	EXPRESSION TAG	UNP Q9H074
A	77	SER	-	EXPRESSION TAG	UNP Q9H074
B	61	GLY	-	EXPRESSION TAG	UNP Q9H074
B	62	SER	-	EXPRESSION TAG	UNP Q9H074
B	63	HIS	-	EXPRESSION TAG	UNP Q9H074
B	64	MET	-	EXPRESSION TAG	UNP Q9H074
B	65	ALA	-	EXPRESSION TAG	UNP Q9H074
B	66	SER	-	EXPRESSION TAG	UNP Q9H074
B	67	MET	-	EXPRESSION TAG	UNP Q9H074
B	68	THR	-	EXPRESSION TAG	UNP Q9H074

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Chain	Residue	Modelled	Actual	Comment	Reference
B	69	GLY	-	EXPRESSION TAG	UNP Q9H074
B	70	GLY	-	EXPRESSION TAG	UNP Q9H074
B	71	GLN	-	EXPRESSION TAG	UNP Q9H074
B	72	GLN	-	EXPRESSION TAG	UNP Q9H074
B	73	MET	-	EXPRESSION TAG	UNP Q9H074
B	74	GLY	-	EXPRESSION TAG	UNP Q9H074
B	75	ARG	-	EXPRESSION TAG	UNP Q9H074
B	76	GLY	-	EXPRESSION TAG	UNP Q9H074
B	77	SER	-	EXPRESSION TAG	UNP Q9H074

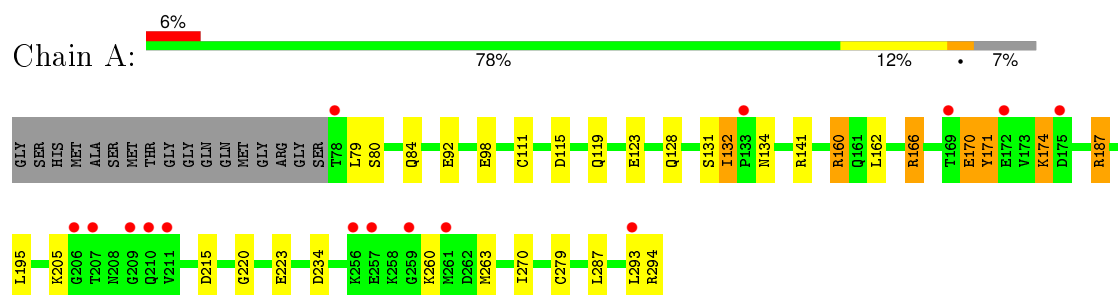
- Molecule 2 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	178	Total O 178 178	0	0
2	B	192	Total O 192 192	0	0

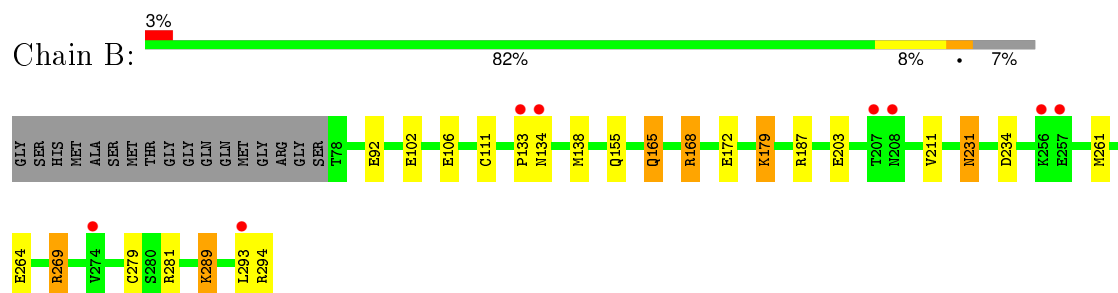
3 Residue-property plots [i](#)

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 ($\text{RSRZ} > 2$). 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: Polyadenylate-binding protein-interacting protein 1



- Molecule 1: Polyadenylate-binding protein-interacting protein 1



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	58.61Å 76.05Å 62.42Å 90.00° 96.47° 90.00°	Depositor
Resolution (Å)	19.95 – 2.00 19.95 – 2.00	Depositor EDS
% Data completeness (in resolution range)	99.5 (19.95-2.00) 99.5 (19.95-2.00)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	0.03	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.25 (at 2.01Å)	Xtriage
Refinement program	REFMAC 5.5.0109	Depositor
R, R_{free}	0.190 , 0.244 0.190 , 0.247	Depositor DCC
R_{free} test set	1834 reflections (5.26%)	DCC
Wilson B-factor (Å ²)	24.6	Xtriage
Anisotropy	0.182	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.40 , 54.0	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.55$, $\langle L^2 \rangle = 0.39$	Xtriage
Outliers	2 of 36673 reflections (0.005%)	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	3952	wwPDB-VP
Average B, all atoms (Å ²)	27.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 40.42 % of the origin peak, indicating pseudo translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo translational symmetry is equal to 2.7433e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.375 respectively for untwinned datasets, and 0.333, 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	0.94	5/1832 (0.3%)	0.85	7/2473 (0.3%)
1	B	1.00	2/1852 (0.1%)	0.89	8/2499 (0.3%)
All	All	0.97	7/3684 (0.2%)	0.87	15/4972 (0.3%)

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	160[A]	ARG	N-CA	-5.75	1.34	1.46
1	A	160[B]	ARG	N-CA	-5.75	1.34	1.46
1	A	171	TYR	CD1-CE1	-5.63	1.30	1.39
1	B	165[A]	GLN	N-CA	-5.27	1.35	1.46
1	B	165[B]	GLN	N-CA	-5.27	1.35	1.46
1	A	160[A]	ARG	CA-C	5.25	1.66	1.52
1	A	160[B]	ARG	CA-C	5.25	1.66	1.52

All (15) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	160[A]	ARG	NE-CZ-NH1	6.84	123.72	120.30
1	A	160[B]	ARG	NE-CZ-NH1	6.84	123.72	120.30
1	B	269	ARG	NE-CZ-NH1	-6.18	117.21	120.30
1	A	195	LEU	CA-CB-CG	6.13	129.41	115.30
1	B	172[A]	GLU	CA-C-N	5.84	130.06	117.20
1	B	172[B]	GLU	CA-C-N	5.84	130.06	117.20
1	B	172[A]	GLU	CA-C-O	-5.80	107.92	120.10
1	B	172[B]	GLU	CA-C-O	-5.80	107.92	120.10
1	A	160[A]	ARG	NE-CZ-NH2	-5.76	117.42	120.30
1	A	160[B]	ARG	NE-CZ-NH2	-5.76	117.42	120.30
1	B	231[A]	ASN	N-CA-CB	-5.48	100.73	110.60
1	B	231[B]	ASN	N-CA-CB	-5.48	100.73	110.60
1	B	281	ARG	NE-CZ-NH1	5.28	122.94	120.30
1	A	160[A]	ARG	CA-C-O	5.15	130.92	120.10
1	A	160[B]	ARG	CA-C-O	5.15	130.92	120.10

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-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 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	1784	0	1796	26	0
1	B	1798	0	1819	39	0
2	A	178	0	0	12	0
2	B	192	0	0	16	0
All	All	3952	0	3615	64	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 9.

All (64) 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:187:ARG:HD2	2:A:314:HOH:O	1.63	0.99
1:B:179[A]:LYS:CB	1:B:179[A]:LYS:NZ	2.27	0.98
1:A:111:CYS:HB3	2:A:364:HOH:O	1.63	0.98
1:B:179[A]:LYS:HB3	1:B:179[A]:LYS:NZ	1.82	0.94
1:B:179[A]:LYS:HZ2	1:B:179[A]:LYS:HB2	1.34	0.91
1:B:179[A]:LYS:HB3	1:B:179[A]:LYS:HZ3	1.35	0.87
1:B:179[A]:LYS:CB	1:B:179[A]:LYS:HZ2	1.86	0.87
1:B:264[B]:GLU:H	1:B:264[B]:GLU:CD	1.85	0.80
1:A:111:CYS:SG	2:A:364:HOH:O	2.42	0.77
1:B:261:MET:O	1:B:264[B]:GLU:OE1	2.02	0.77
1:A:92:GLU:OE1	2:A:445:HOH:O	2.04	0.76
1:B:168[A]:ARG:NH2	2:B:357:HOH:O	1.81	0.74
1:B:111:CYS:HB3	2:B:323:HOH:O	1.88	0.73
1:B:111:CYS:SG	2:B:323:HOH:O	2.46	0.73
1:B:138:MET:HE1	2:B:324:HOH:O	1.87	0.73
1:A:123:GLU:OE1	1:A:166:ARG:NH1	2.25	0.69
1:B:269:ARG:NH1	2:B:384:HOH:O	2.19	0.66
1:A:160[A]:ARG:NH2	2:A:444:HOH:O	2.31	0.62
1:B:261:MET:C	1:B:264[B]:GLU:OE1	2.38	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:133:PRO:HA	2:B:327:HOH:O	2.01	0.60
1:A:84:GLN:HE22	1:A:128:GLN:NE2	1.99	0.59
1:B:92:GLU:OE1	2:B:439:HOH:O	2.17	0.58
1:A:111:CYS:CB	2:A:364:HOH:O	2.28	0.58
1:B:155:GLN:HG2	2:B:442:HOH:O	2.05	0.57
1:A:171:TYR:CE1	1:A:220:GLY:HA2	2.41	0.56
1:B:187:ARG:NH2	2:B:412:HOH:O	2.21	0.56
1:A:84:GLN:HE22	1:A:128:GLN:HE22	1.54	0.56
1:B:261:MET:HE1	2:B:360:HOH:O	2.07	0.54
1:A:119[A]:GLN:HG3	1:A:162:LEU:HD13	1.89	0.54
1:A:294:ARG:CG	2:A:347:HOH:O	2.57	0.52
1:B:134[A]:ASN:ND2	2:B:422:HOH:O	2.42	0.50
1:A:234:ASP:OD1	1:A:279:CYS:HB2	2.11	0.50
1:A:170[B]:GLU:OE2	2:A:352:HOH:O	2.19	0.50
1:B:203:GLU:HB3	1:B:211:VAL:HG13	1.93	0.50
1:A:187:ARG:CD	2:A:314:HOH:O	2.40	0.50
1:B:293:LEU:CD1	2:B:411:HOH:O	2.60	0.49
1:A:293:LEU:HD12	2:A:418:HOH:O	2.11	0.49
1:B:133:PRO:O	1:B:134[B]:ASN:CG	2.52	0.48
1:B:102:GLU:O	1:B:106[B]:GLU:HG3	2.13	0.48
1:A:98:GLU:CD	1:A:141:ARG:HH22	2.18	0.47
1:A:171:TYR:OH	1:A:223:GLU:HB2	2.16	0.46
1:A:294:ARG:HG2	2:A:347:HOH:O	2.15	0.46
1:B:165[B]:GLN:HE22	1:B:168[B]:ARG:HE	1.63	0.46
1:A:134:ASN:ND2	1:B:134[A]:ASN:OD1	2.41	0.45
1:B:138:MET:CE	2:B:324:HOH:O	2.56	0.45
1:B:289:LYS:O	1:B:293:LEU:HD13	2.17	0.45
1:A:270:ILE:HG23	1:A:287:LEU:HB3	1.99	0.45
1:B:179[B]:LYS:HD2	2:B:404:HOH:O	2.18	0.44
1:B:165[B]:GLN:NE2	1:B:168[B]:ARG:HE	2.17	0.43
1:A:263:MET:HE1	1:A:294:ARG:HG3	1.99	0.43
1:B:234:ASP:OD1	1:B:279:CYS:HB2	2.18	0.43
1:B:179[A]:LYS:NZ	1:B:231[A]:ASN:OD1	2.49	0.43
1:B:293:LEU:HD11	2:B:411:HOH:O	2.19	0.42
1:A:174:LYS:HE2	1:A:223:GLU:HB3	2.01	0.42
1:A:187:ARG:NH1	2:A:328:HOH:O	2.50	0.42
1:B:155:GLN:NE2	2:B:333:HOH:O	2.51	0.41
1:B:102:GLU:O	1:B:106[B]:GLU:CG	2.68	0.41
1:A:171:TYR:OH	1:A:223:GLU:HG3	2.21	0.41
1:A:131:SER:O	1:A:132:ILE:C	2.60	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	223/234 (95%)	218 (98%)	4 (2%)	1 (0%)	39	33
1	B	225/234 (96%)	224 (100%)	1 (0%)	0	100	100
All	All	448/468 (96%)	442 (99%)	5 (1%)	1 (0%)	52	48

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	132	ILE

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	202/205 (98%)	191 (95%)	11 (5%)	27	21
1	B	204/205 (100%)	198 (97%)	6 (3%)	50	49
All	All	406/410 (99%)	389 (96%)	17 (4%)	42	31

All (17) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	79	LEU
1	A	80	SER
1	A	115	ASP
1	A	166	ARG
1	A	170[A]	GLU

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Mol	Chain	Res	Type
1	A	170[B]	GLU
1	A	174	LYS
1	A	187	ARG
1	A	205	LYS
1	A	215	ASP
1	A	260	LYS
1	B	168[A]	ARG
1	B	168[B]	ARG
1	B	179[A]	LYS
1	B	179[B]	LYS
1	B	289	LYS
1	B	294	ARG

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

Mol	Chain	Res	Type
1	A	128	GLN
1	A	134	ASN
1	A	148	HIS
1	A	201	ASN
1	A	218	GLN
1	B	148	HIS
1	B	236	ASN

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 carbohydrates in this entry.

5.6 Ligand geometry [i](#)

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	A	217/234 (92%)	0.18	15 (6%) 20 21	14, 26, 46, 53	0
1	B	217/234 (92%)	0.00	8 (3%) 45 47	13, 23, 40, 52	0
All	All	434/468 (92%)	0.09	23 (5%) 30 32	13, 25, 44, 53	0

All (23) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	207	THR	6.2
1	B	133	PRO	4.4
1	A	293	LEU	3.9
1	B	293	LEU	3.9
1	A	206	GLY	3.5
1	A	172	GLU	3.4
1	A	211	VAL	3.2
1	A	261	MET	3.1
1	A	207	THR	3.1
1	A	209	GLY	3.0
1	A	78	THR	3.0
1	A	133	PRO	2.8
1	B	257	GLU	2.7
1	A	175	ASP	2.6
1	A	256	LYS	2.6
1	B	256	LYS	2.5
1	B	208	ASN	2.4
1	B	134[A]	ASN	2.4
1	A	169	THR	2.3
1	A	210	GLN	2.3
1	A	257	GLU	2.2
1	A	259	GLY	2.1
1	B	274	VAL	2.1

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 carbohydrates 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.